

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
 Data File : BM047506.D
 Acq On : 12 Sep 2024 16:49
 Operator : RC/JU
 Sample : P3878-08DL2 20X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVY6DL2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M

Title : SVOA CALIBRATION

Signal : TIC: BM047506.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.999	131	155	158	rBV	690649	2485655	3.21%	0.732%
2	4.975	316	321	326	rVV	613889	872489	1.13%	0.257%
3	5.899	471	478	487	rVB2	2312859	3700159	4.77%	1.090%
4	6.175	514	525	531	rBV	1487677	2875480	3.71%	0.847%
5	6.369	552	558	565	rVB	971149	1550685	2.00%	0.457%
6	6.440	565	570	575	rBV	864479	1233485	1.59%	0.363%
7	6.516	575	583	586	rBV3	634835	1342735	1.73%	0.396%
8	6.551	586	589	595	rVB	590781	878358	1.13%	0.259%
9	6.781	619	628	633	rVB4	385649	1024621	1.32%	0.302%
10	6.846	633	639	640	rBV	661360	990243	1.28%	0.292%
11	6.875	640	644	649	rVB	1226165	1891573	2.44%	0.557%
12	6.934	649	654	657	rBV	1544315	2433934	3.14%	0.717%
13	6.963	657	659	664	rVB	1202065	1559758	2.01%	0.460%
14	7.040	664	672	677	rBV2	1679131	3838702	4.95%	1.131%
15	7.098	677	682	687	rVB	796597	1188415	1.53%	0.350%
16	7.263	702	710	713	rBV2	898168	1508732	1.95%	0.445%
17	7.298	713	716	720	rVB	1173712	1460866	1.88%	0.430%
18	7.393	725	732	737	rVB2	1084409	2003726	2.58%	0.590%
19	7.510	746	752	759	rVB2	9348473	15238482	19.65%	4.490%
20	7.716	779	787	791	rBV3	342605	893548	1.15%	0.263%
21	7.863	806	812	818	rVB	3273148	5101567	6.58%	1.503%
22	7.934	818	824	830	rBV	4323493	7247107	9.35%	2.135%
23	7.987	830	833	843	rVB2	779845	1444278	1.86%	0.426%
24	8.098	843	852	856	rBV2	1569466	3240390	4.18%	0.955%
25	8.169	861	864	871	rVB3	749249	1275215	1.64%	0.376%
26	8.298	881	886	891	rBV	1066030	1988701	2.56%	0.586%
27	8.416	901	906	912	rVB2	1636176	2743584	3.54%	0.808%
28	8.475	912	916	919	rBV	997769	1236172	1.59%	0.364%
29	8.516	919	923	925	rBV2	1258286	1745104	2.25%	0.514%
30	8.545	925	928	937	rVB	1844705	3213014	4.14%	0.947%
31	8.651	940	946	949	rBV	1499076	2560567	3.30%	0.754%
32	8.828	971	976	980	rBV	761758	1080268	1.39%	0.318%
33	8.875	980	984	993	rVB	805889	1355428	1.75%	0.399%
34	8.981	993	1002	1012	rBV4	1183971	3141800	4.05%	0.926%
35	9.110	1013	1024	1029	rBV	7326628	11281248	14.55%	3.324%
36	9.234	1036	1045	1051	rBV8	856528	2444612	3.15%	0.720%

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Integrator: RTE

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Title : SVOA CALIBRATION

37	9.310	1051	1058	1062	rBV3	540174	1149415	1.48%	0.339%
38	9.675	1112	1120	1125	rVB2	726594	1340542	1.73%	0.395%
39	9.816	1139	1144	1147	rBV4	994766	1924971	2.48%	0.567%
40	9.963	1160	1169	1174	rVB	950280	2037365	2.63%	0.600%
41	10.028	1174	1180	1184	rBV2	679986	1123784	1.45%	0.331%
42	10.110	1190	1194	1197	rVV	715760	1152832	1.49%	0.340%
43	10.216	1207	1212	1216	rBV2	515081	856389	1.10%	0.252%
44	10.369	1232	1238	1243	rBV2	854675	1377545	1.78%	0.406%
45	10.545	1262	1268	1272	rBV4	418794	835293	1.08%	0.246%
46	10.657	1280	1287	1294	rVV2	6708567	13091076	16.88%	3.857%
47	10.722	1295	1298	1305	rVV2	721992	1317714	1.70%	0.388%
48	10.792	1305	1310	1315	rVV2	1288611	2285268	2.95%	0.673%
49	10.845	1315	1319	1324	rVV	599998	916088	1.18%	0.270%
50	11.057	1348	1355	1367	rBV	3542940	5811818	7.50%	1.713%
51	11.175	1367	1375	1382	rBV2	1365840	2762492	3.56%	0.814%
52	11.598	1437	1447	1459	rVB9	469526	1648834	2.13%	0.486%
53	11.704	1459	1465	1470	rBV2	1470871	2717879	3.51%	0.801%
54	11.757	1470	1474	1484	rVB3	618556	1462675	1.89%	0.431%
55	11.939	1498	1505	1514	rBV	2027460	3318288	4.28%	0.978%
56	12.104	1525	1533	1535	rBV	2459156	3884421	5.01%	1.145%
57	12.133	1535	1538	1544	rVB	2037408	2934634	3.78%	0.865%
58	12.339	1565	1573	1579	rVB2	3835888	7113528	9.17%	2.096%
59	12.504	1596	1601	1604	rBV4	630810	1002357	1.29%	0.295%
60	12.545	1604	1608	1612	rVV2	546116	869759	1.12%	0.256%
61	12.916	1666	1671	1680	rVB2	611456	1246754	1.61%	0.367%
62	13.057	1691	1695	1702	rVB2	860848	1290771	1.66%	0.380%
63	13.339	1737	1743	1750	rBV	2918109	4628375	5.97%	1.364%
64	13.469	1761	1765	1773	rVV2	453131	857720	1.11%	0.253%
65	13.663	1789	1798	1799	rBV3	759047	1320585	1.70%	0.389%
66	13.769	1812	1816	1820	rBV	956395	1286805	1.66%	0.379%
67	13.816	1820	1824	1828	rBV	786639	1355226	1.75%	0.399%
68	13.998	1846	1855	1859	rBV2	1279738	2389858	3.08%	0.704%
69	14.045	1859	1863	1868	rVV4	578060	1093202	1.41%	0.322%
70	14.410	1921	1925	1930	rBV	3731307	4752844	6.13%	1.400%
71	14.492	1934	1939	1945	rVB	4242046	6563242	8.46%	1.934%
72	14.580	1950	1954	1960	rVB4	971010	1427259	1.84%	0.421%
73	14.798	1985	1991	1996	rBV4	444099	1025414	1.32%	0.302%
74	14.969	2014	2020	2024	rVB	677896	1024526	1.32%	0.302%
75	15.033	2026	2031	2035	rVV3	972717	1529328	1.97%	0.451%
76	15.116	2039	2045	2049	rVV	5258237	8050675	10.38%	2.372%

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Title : SVOA CALIBRATION

77	15.157	2049	2052	2057	rVB3	654286	974983	1.26%	0.287%
78	15.363	2083	2087	2091	rVB	3135594	4033395	5.20%	1.188%
79	15.592	2121	2126	2131	rVB5	531339	907703	1.17%	0.267%
80	15.769	2149	2156	2161	rVB	1109003	2102547	2.71%	0.620%
81	15.980	2188	2192	2200	rVB3	495923	1027520	1.33%	0.303%
82	16.221	2228	2233	2236	rBV	3319785	4944541	6.38%	1.457%
83	16.245	2236	2237	2242	rVB	1494366	1572594	2.03%	0.463%
84	16.892	2339	2347	2352	rBV	42365807	77535208	100.00%	22.846%
85	17.010	2363	2367	2371	rVB	2656696	3114287	4.02%	0.918%
86	17.063	2371	2376	2385	rVB2	1324742	2308652	2.98%	0.680%
87	17.233	2400	2405	2410	rBV	4504493	6169712	7.96%	1.818%
88	17.745	2488	2492	2497	rVB	2519694	2946918	3.80%	0.868%
89	17.915	2516	2521	2525	rBV	1285125	1634467	2.11%	0.482%
90	18.433	2605	2609	2615	rVB	1842605	2282446	2.94%	0.673%
91	19.068	2712	2717	2723	rVV2	1741814	3300239	4.26%	0.972%
92	19.215	2738	2742	2750	rVB	630301	882170	1.14%	0.260%
93	19.645	2811	2815	2825	rVB	969545	1264683	1.63%	0.373%
94	20.174	2901	2905	2911	rBV	921596	1104977	1.43%	0.326%
95	21.156	3067	3072	3082	rVB	2315360	3082241	3.98%	0.908%
96	21.456	3117	3123	3132	rVB	4205471	6221896	8.02%	1.833%
97	21.680	3157	3161	3171	rVB	562924	879826	1.13%	0.259%
98	22.139	3234	3239	3253	rVB	941610	1575437	2.03%	0.464%
99	22.262	3255	3260	3270	rVB2	736662	1296327	1.67%	0.382%
100	24.497	3631	3640	3649	rBV	2182907	5363723	6.92%	1.580%

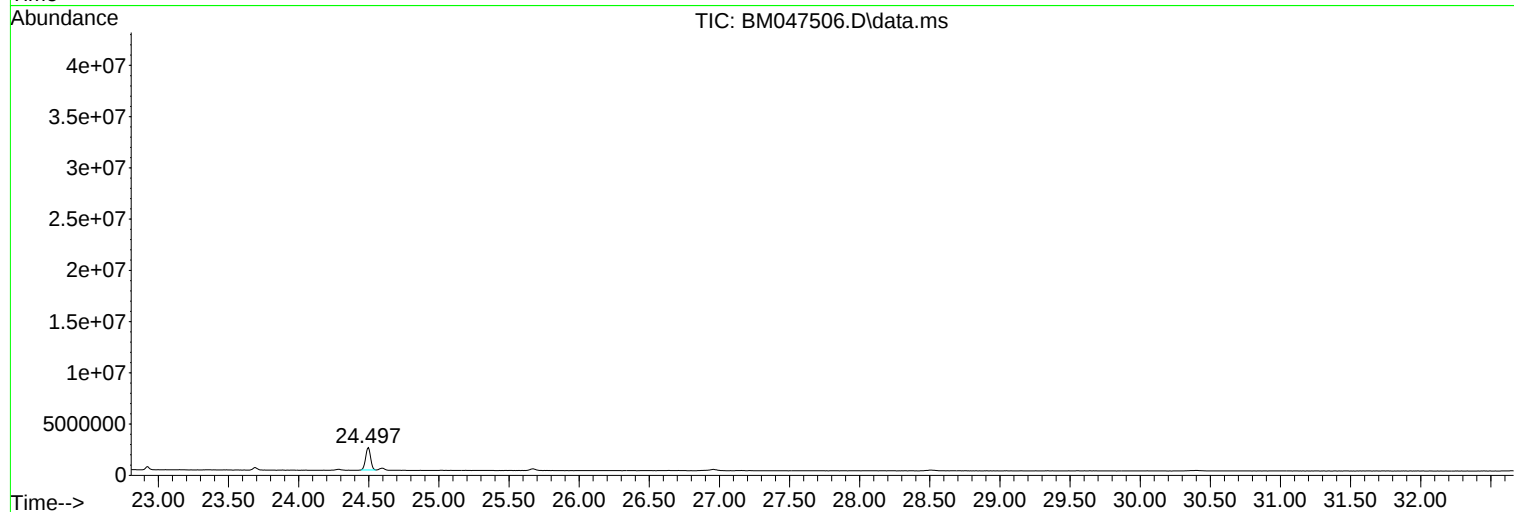
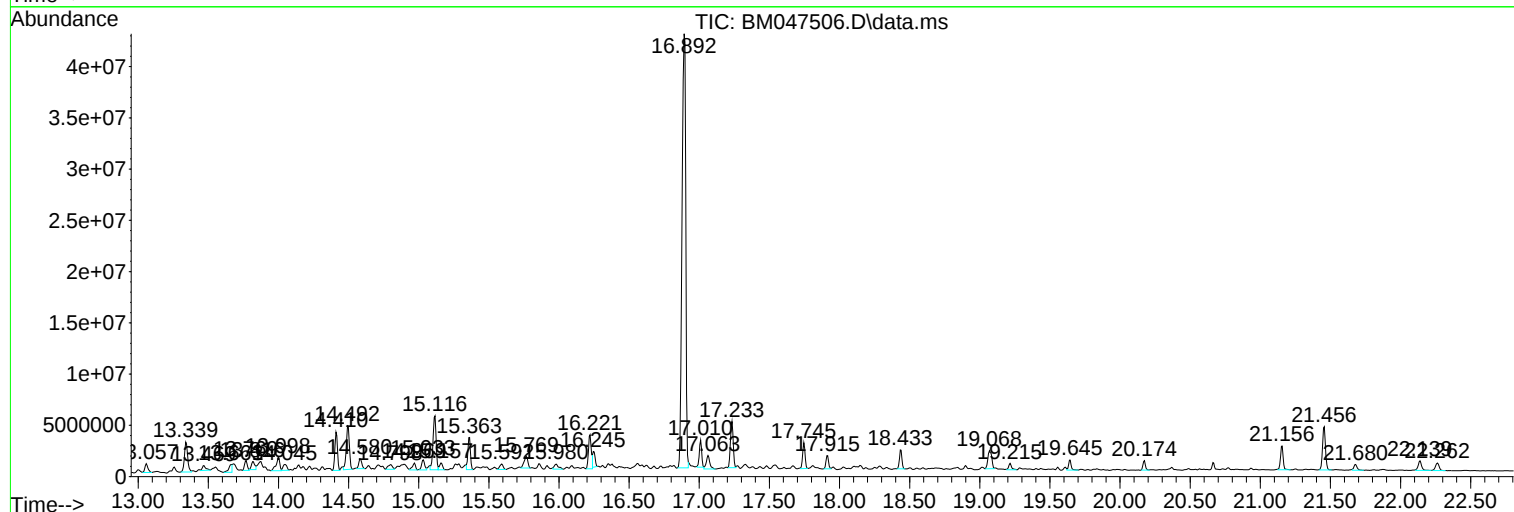
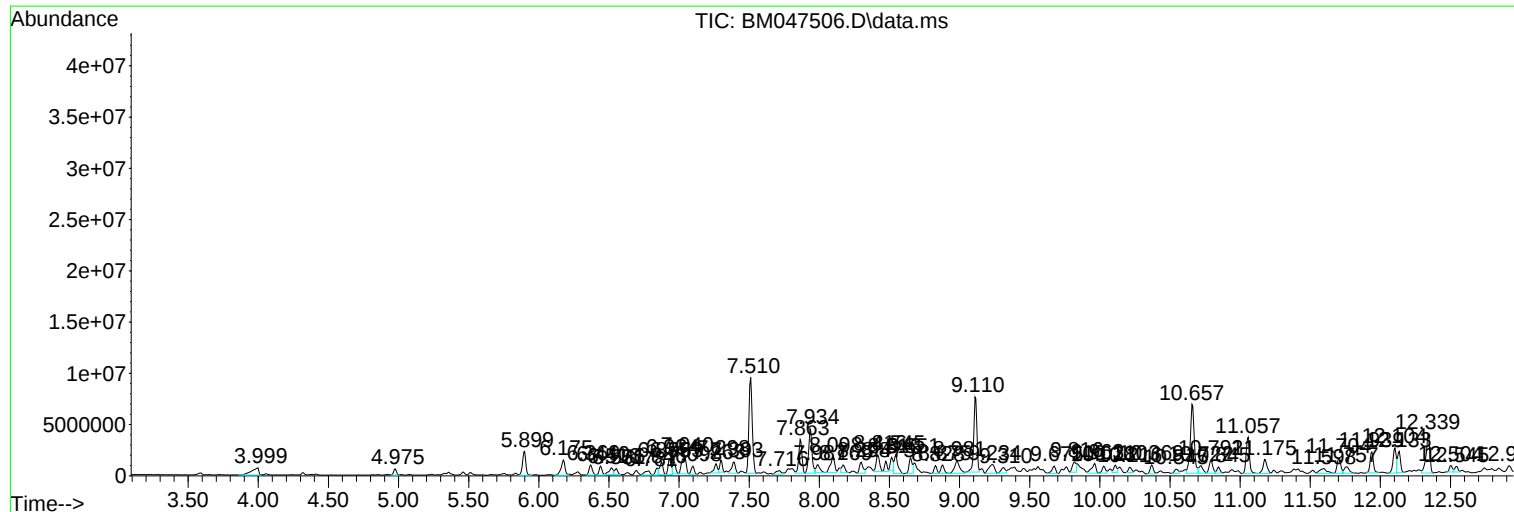
Sum of corrected areas: 339374744

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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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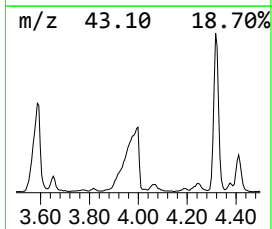
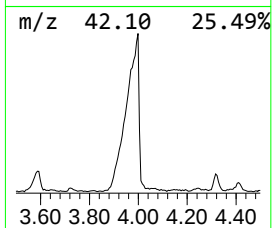
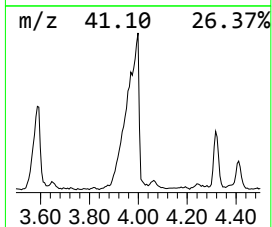
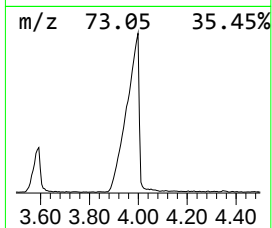
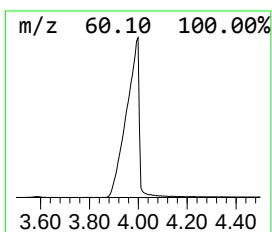
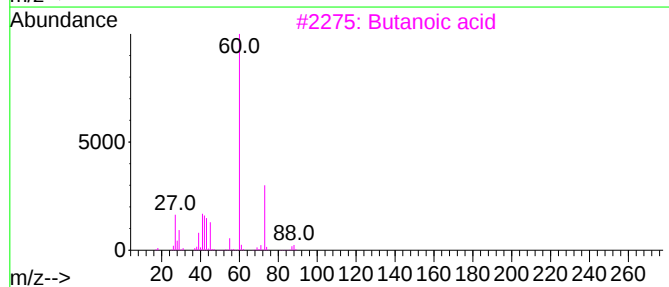
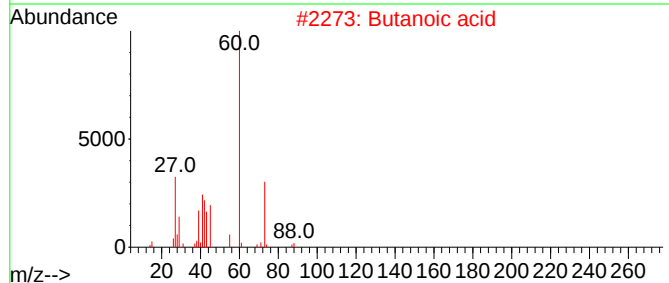
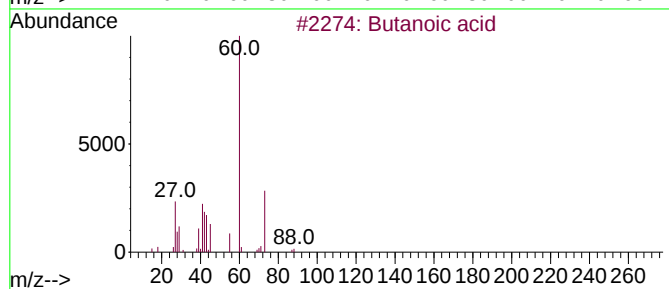
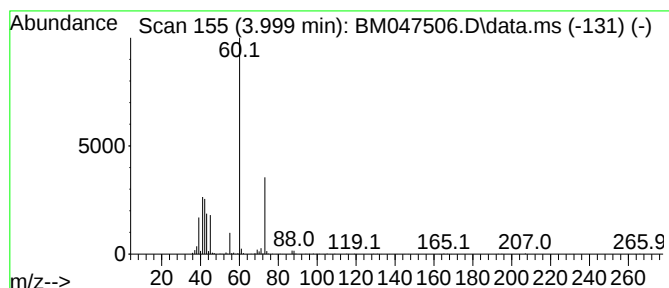
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.999	9.74 ng/ul	2485660	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid	88	C4H8O2	000107-92-6	91
2			Butanoic acid	88	C4H8O2	000107-92-6	90
3			Butanoic acid	88	C4H8O2	000107-92-6	86
4			Butanoic acid	88	C4H8O2	000107-92-6	72
5			Silver butanoate	194	C4H7AgO2	005076-24-4	43



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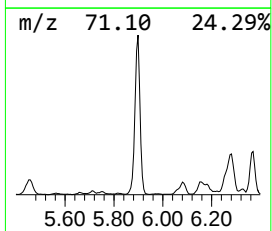
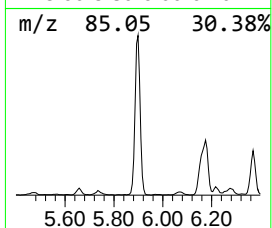
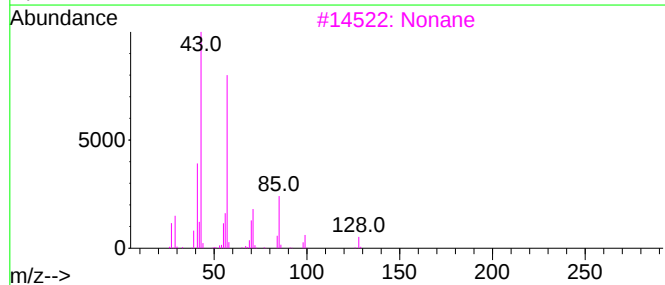
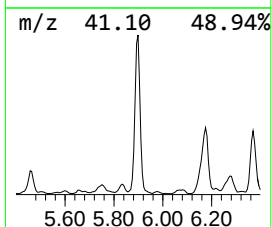
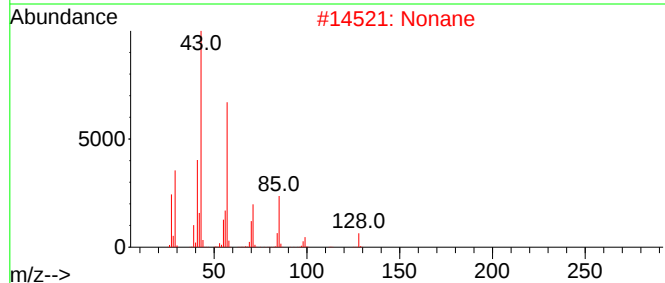
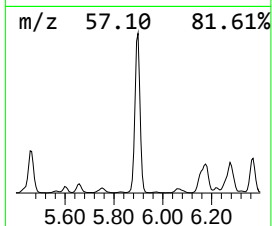
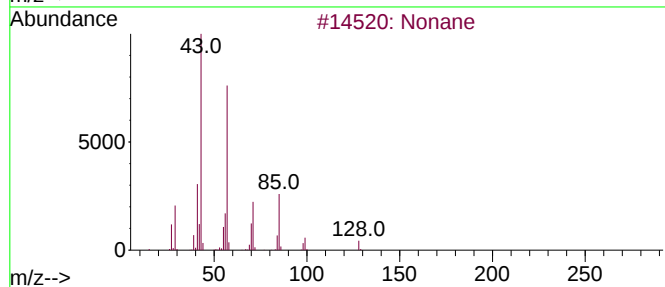
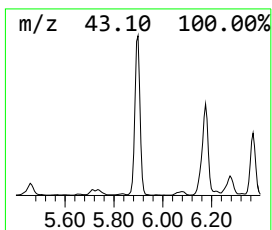
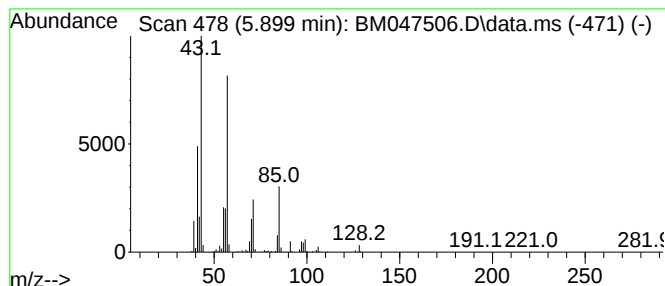
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.899	14.51 ng/ul	3700160	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	91
2	Nonane			128	C9H20	000111-84-2	90
3	Nonane			128	C9H20	000111-84-2	83
4	Octane			114	C8H18	000111-65-9	72
5	Decane			142	C10H22	000124-18-5	64



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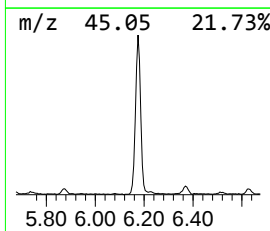
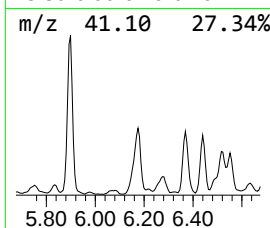
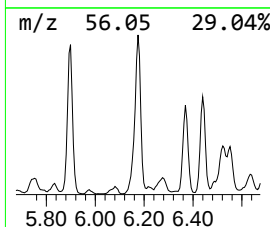
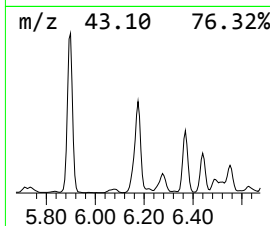
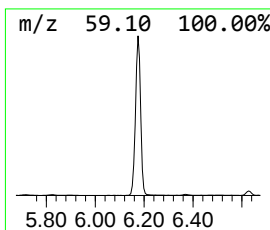
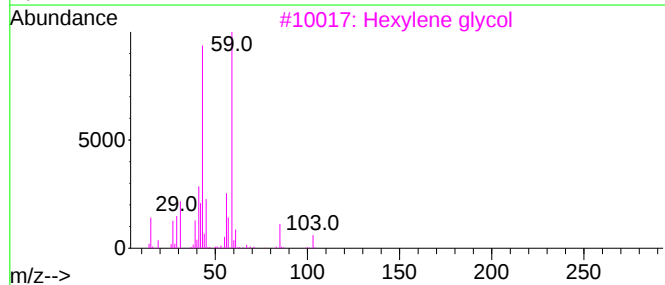
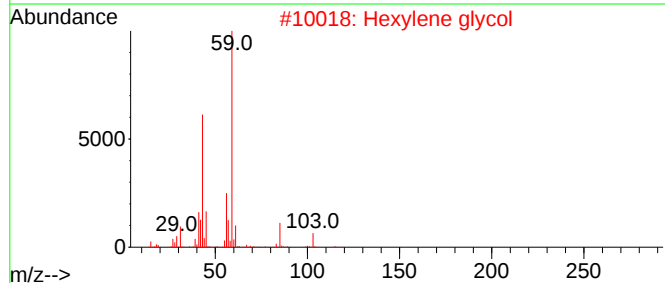
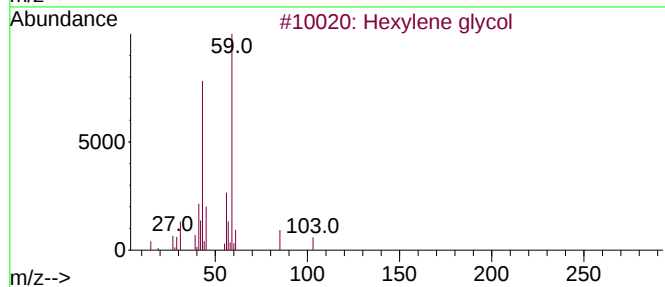
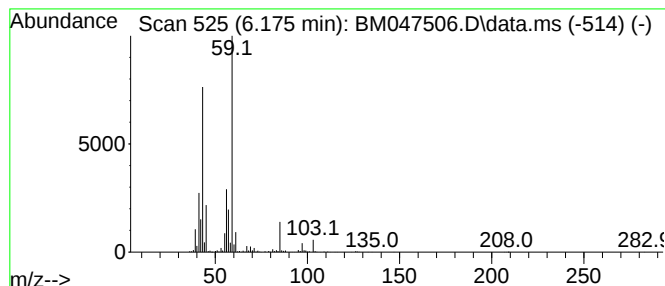
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexylene glycol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.175	11.27 ng/ul	2875480	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	83
3			Hexylene glycol	118	C6H14O2	000107-41-5	83
4			Hexylene glycol	118	C6H14O2	000107-41-5	72
5			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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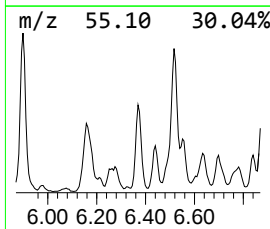
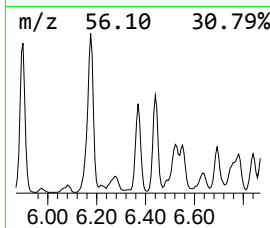
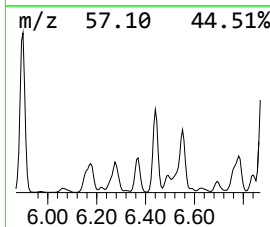
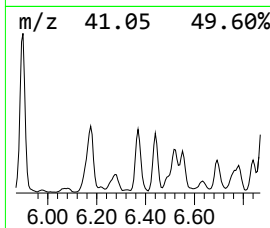
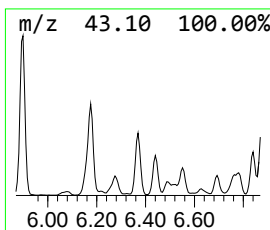
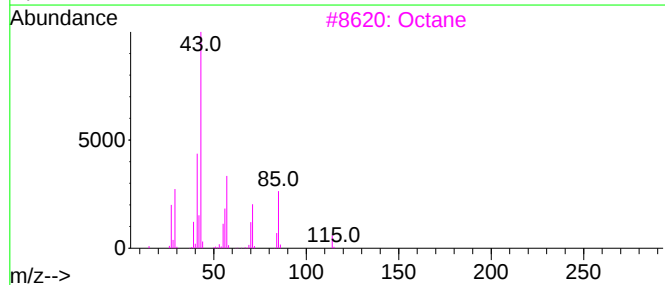
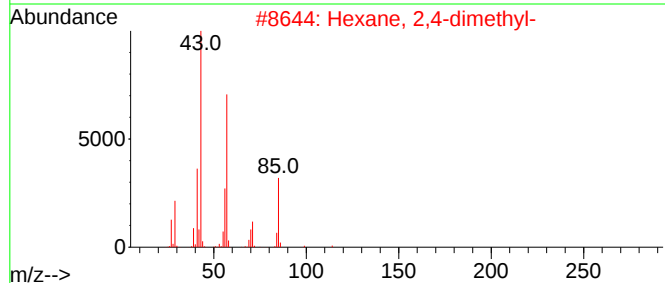
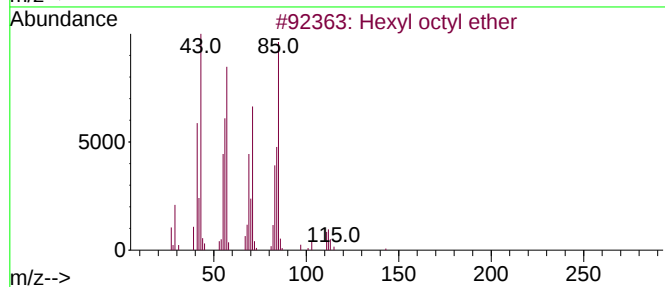
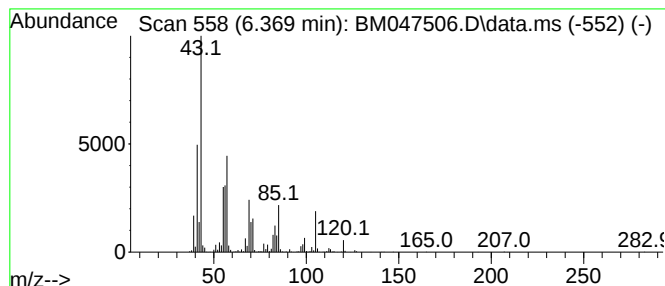
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-01 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.369	6.08 ng/ul	1550690	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexyl octyl ether	214	C14H30O	017071-54-4	47
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	47
3			Octane	114	C8H18	000111-65-9	43
4			Octane	114	C8H18	000111-65-9	43
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
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ClientSampleId :
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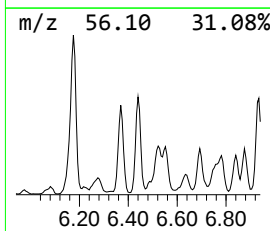
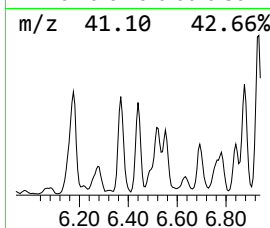
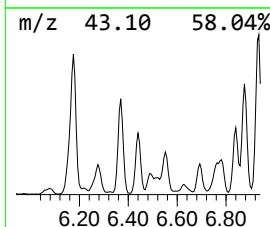
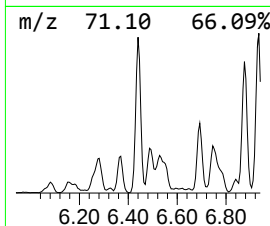
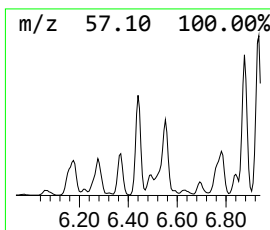
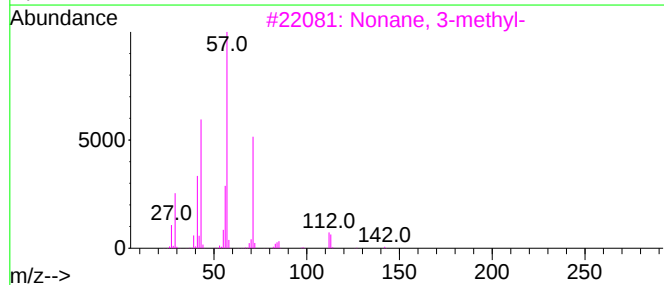
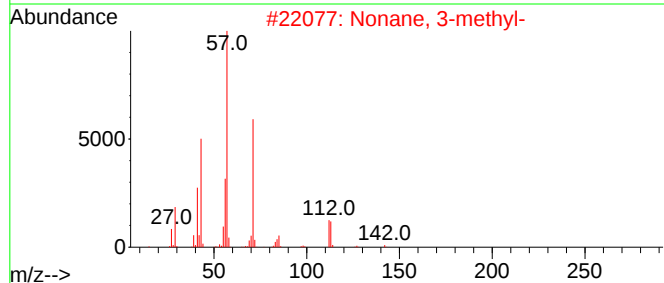
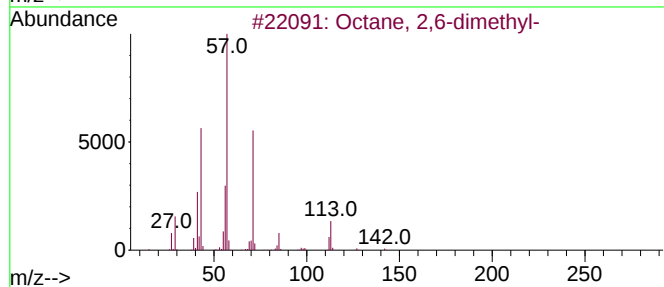
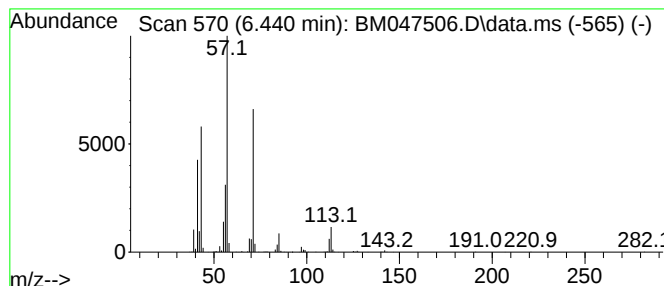
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.440	4.84 ng/ul	1233490	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	93
2	Nonane, 3-methyl-			142	C10H22	005911-04-6	91
3	Nonane, 3-methyl-			142	C10H22	005911-04-6	87
4	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	86
5	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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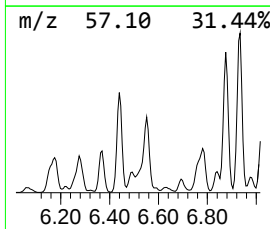
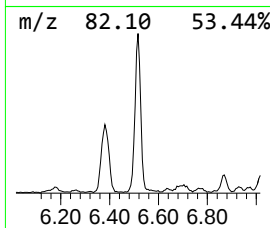
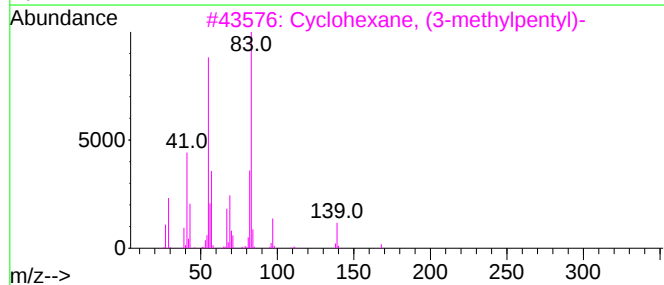
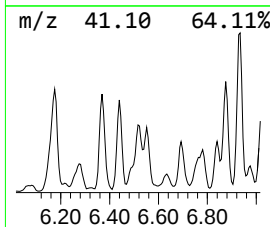
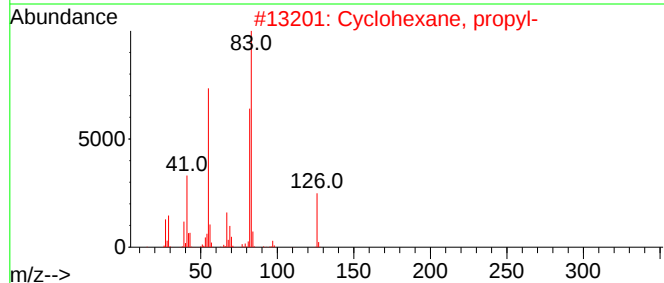
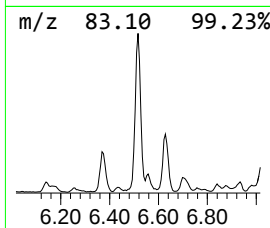
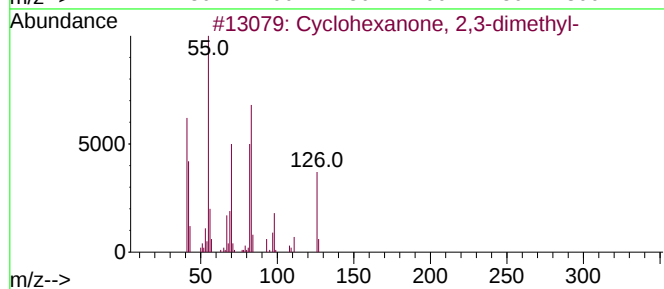
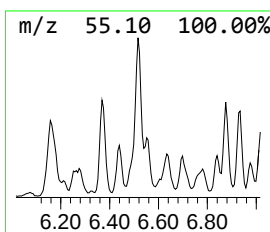
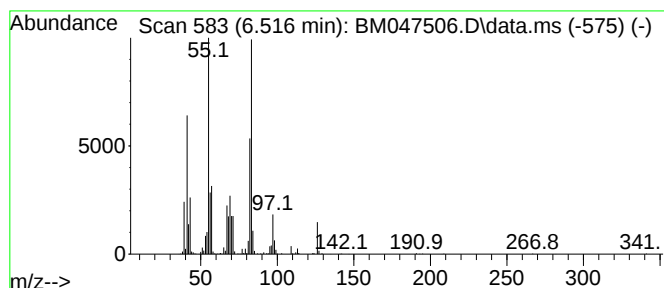
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Cyclohexanone, 2,3-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.516	5.26 ng/ul	1342740	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	87
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	83
3			Cyclohexane, (3-methylpentyl)-	168	C12H24	061142-38-9	78
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	72
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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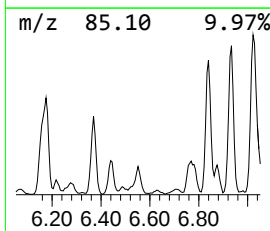
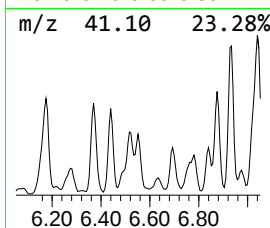
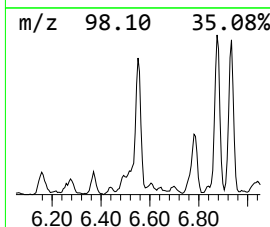
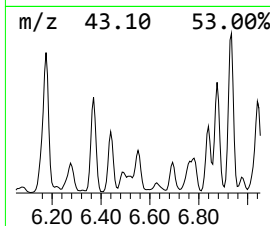
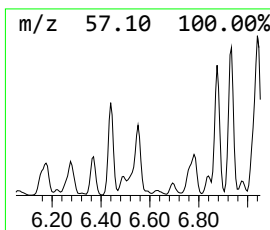
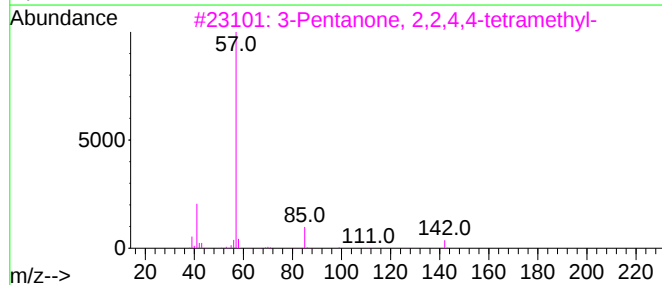
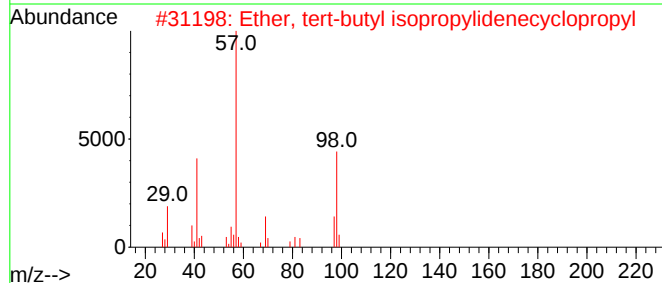
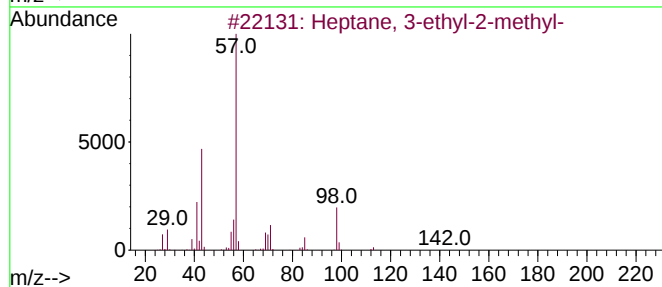
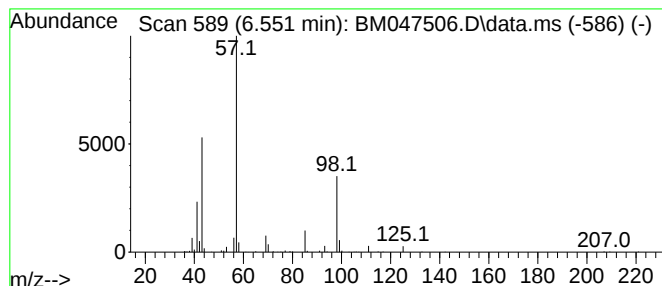
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.551	3.44 ng/ul	878358	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
2			Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	45
3			3-Pentanone, 2,2,4,4-tetramethyl-	142	C9H18O	000815-24-7	38
4			Heptane, 2,3,5-trimethyl-	142	C10H22	020278-85-7	38
5			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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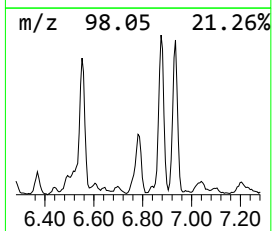
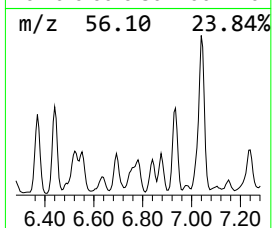
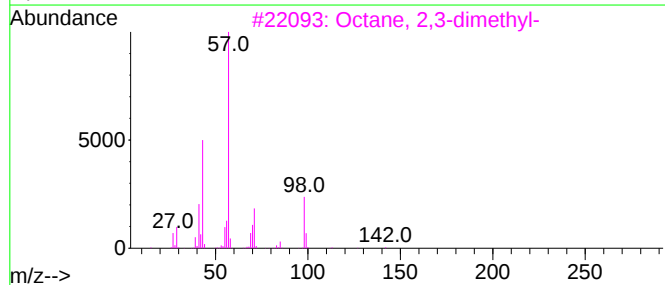
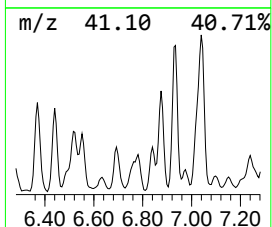
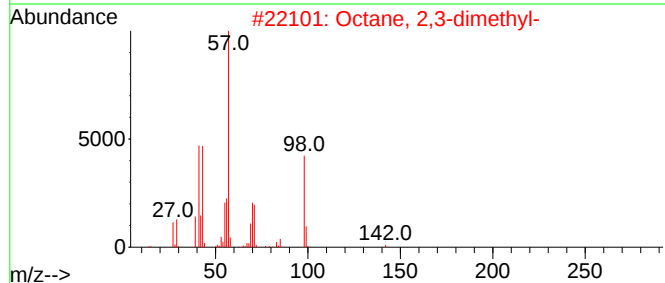
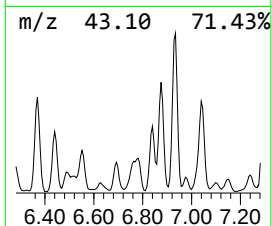
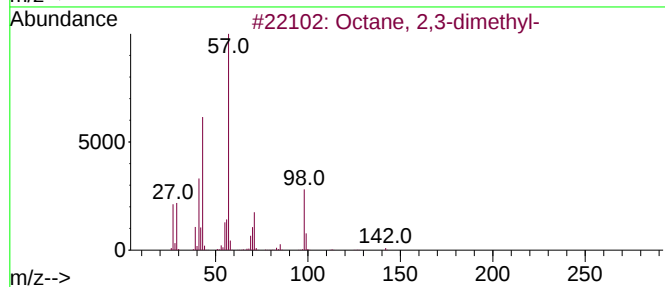
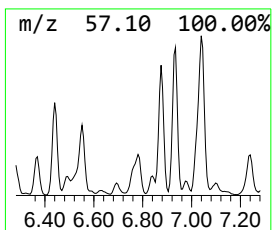
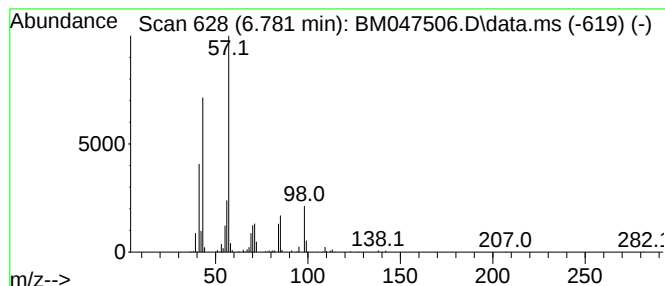
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.781	4.02 ng/ul	1024620	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,3-dimethyl-		142	C10H22	007146-60-3	58	
2	Octane, 2,3-dimethyl-		142	C10H22	007146-60-3	58	
3	Octane, 2,3-dimethyl-		142	C10H22	007146-60-3	53	
4	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	53	
5	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	52	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
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Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
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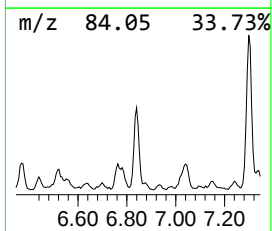
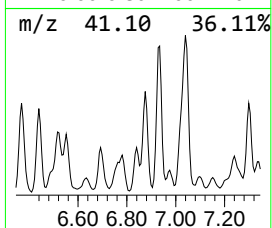
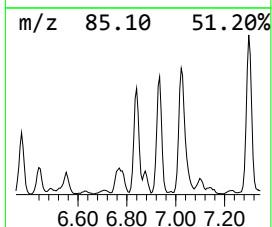
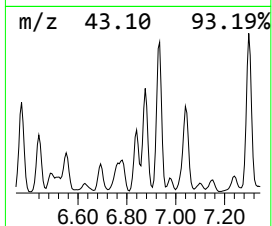
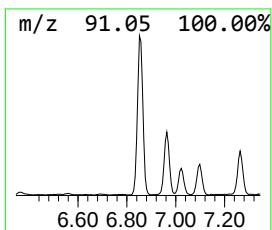
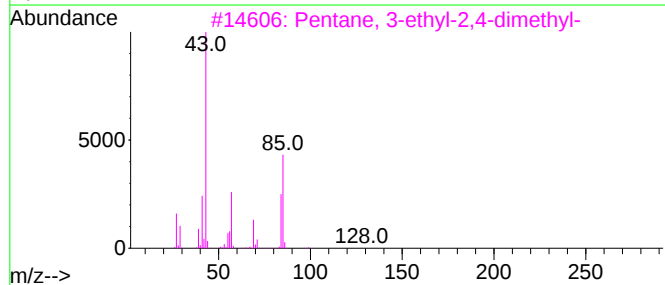
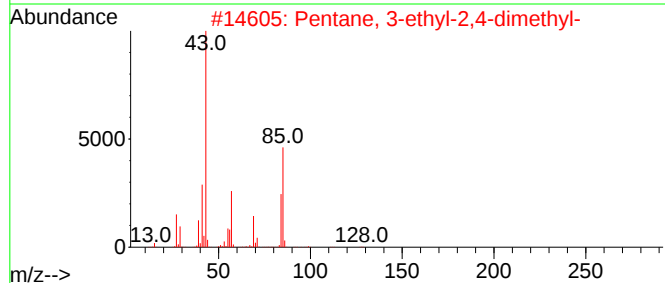
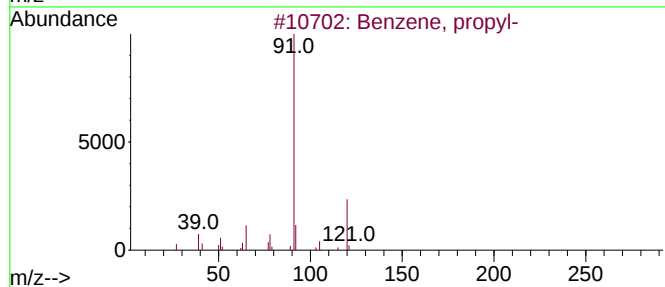
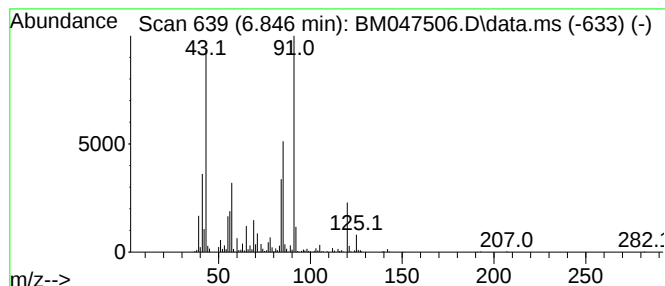
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, propyl- Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.846	3.88 ng/ul	990243	1,4-Dichlorobenzene-d4			7.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, propyl-		120	C9H12	000103-65-1	80
2	Pentane, 3-ethyl-2,4-dimethyl-		128	C9H20	001068-87-7	50
3	Pentane, 3-ethyl-2,4-dimethyl-		128	C9H20	001068-87-7	50
4	Nonane, 5-methyl-		142	C10H22	015869-85-9	46
5	Benzene, propyl-		120	C9H12	000103-65-1	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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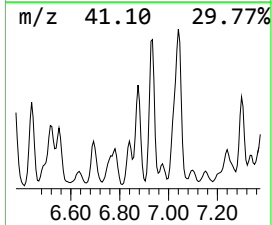
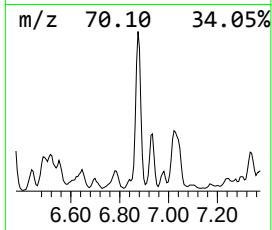
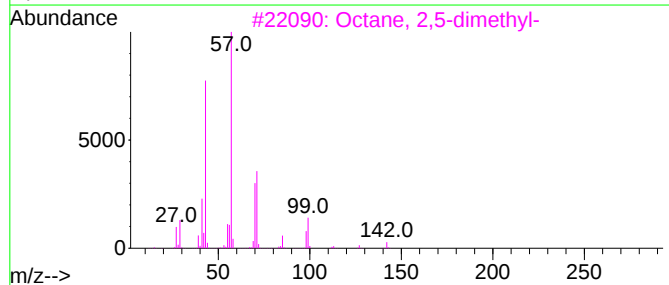
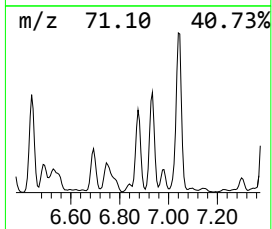
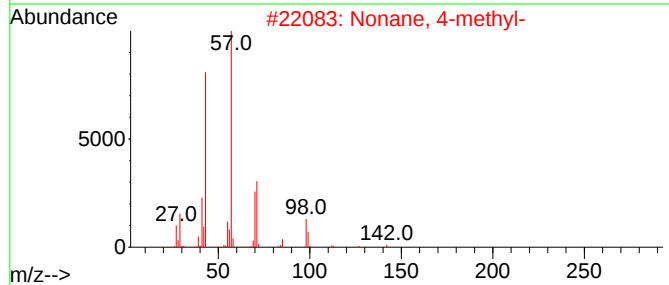
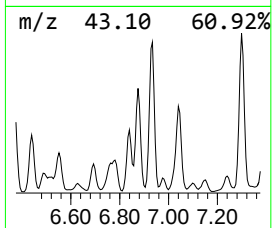
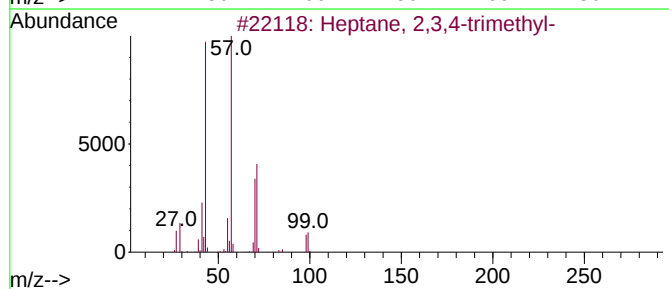
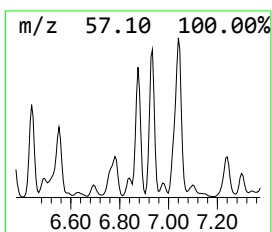
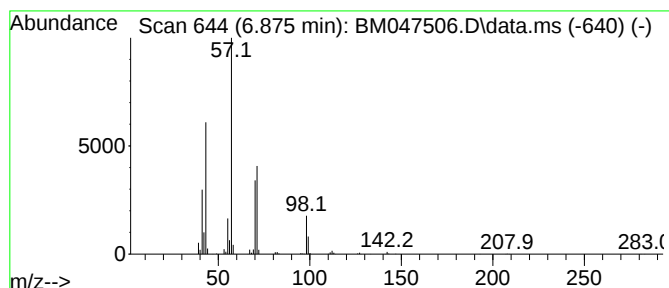
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.875	7.42 ng/ul	1891570	1,4-Dichlorobenzene-d4	7.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	78
2		Nonane, 4-methyl-	142	C10H22	017301-94-9	62
3		Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	56
4		Nonane, 4-methyl-	142	C10H22	017301-94-9	56
5		Nonane, 2-methyl-	142	C10H22	000871-83-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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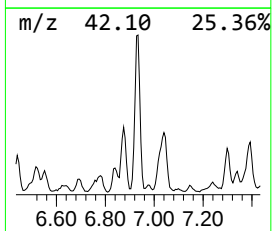
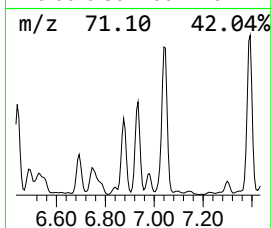
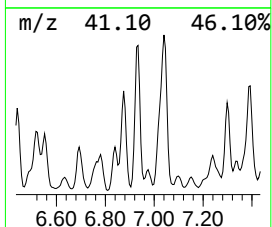
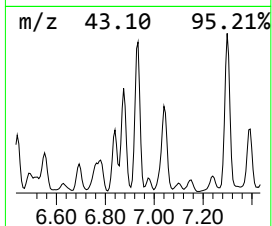
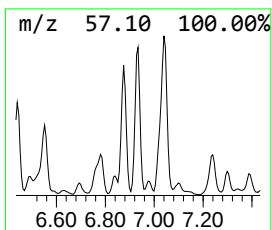
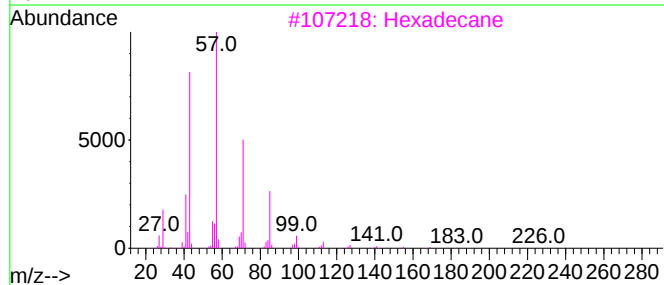
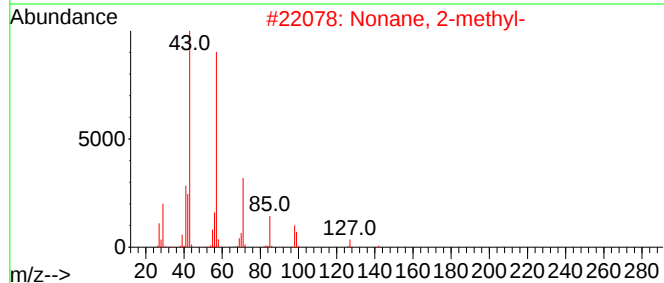
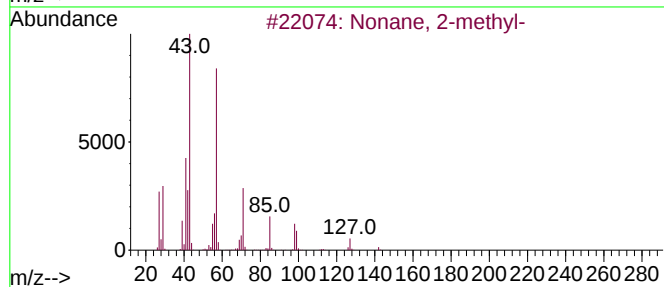
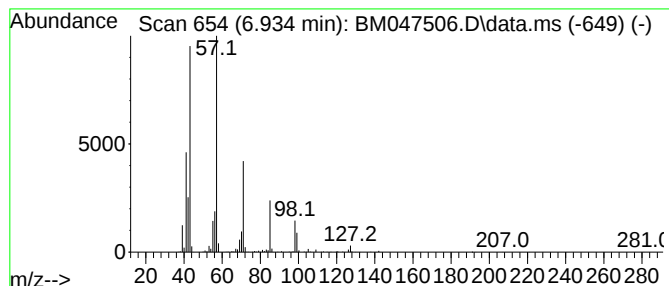
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.934	9.54 ng/ul	2433930	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	83
2			Nonane, 2-methyl-	142	C10H22	000871-83-0	81
3			Hexadecane	226	C16H34	000544-76-3	72
4			Tridecane	184	C13H28	000629-50-5	59
5			Tridecane, 6-methyl-	198	C14H30	013287-21-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

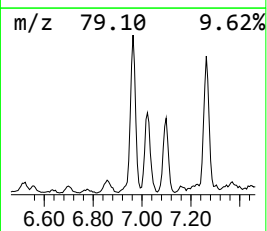
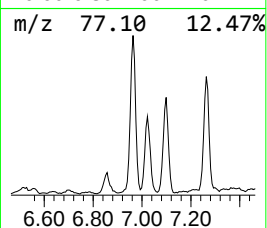
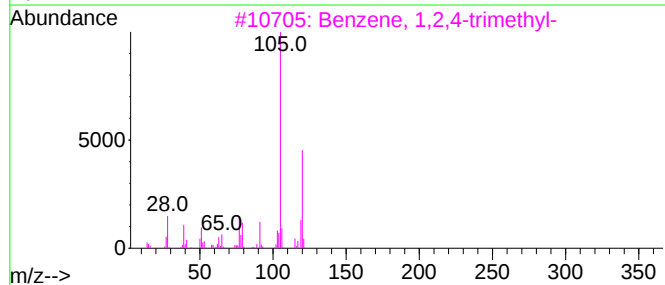
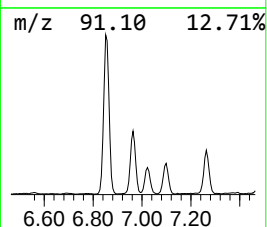
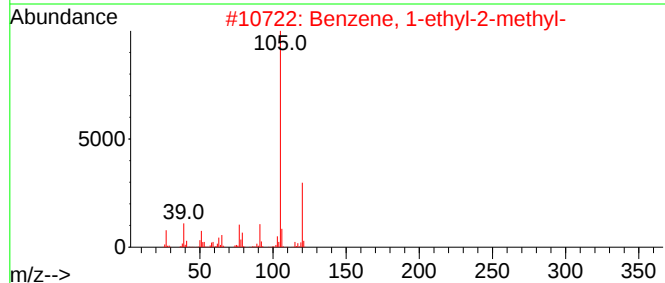
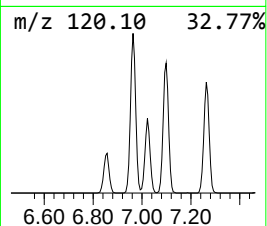
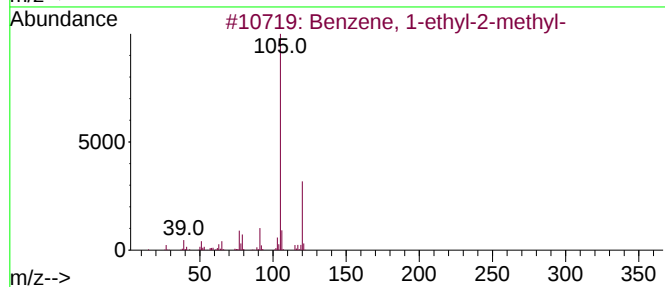
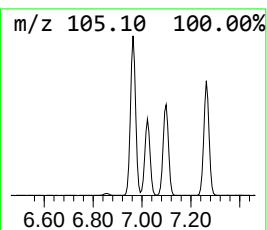
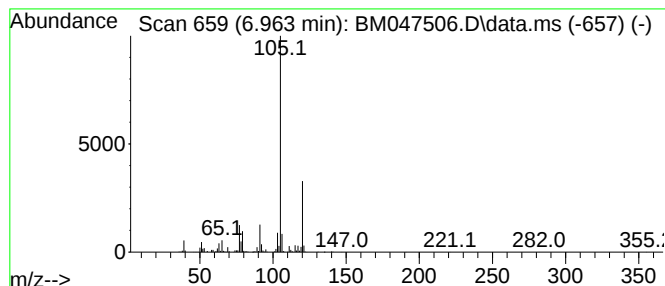
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.963	6.11 ng/ul	1559760	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

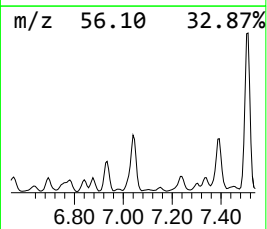
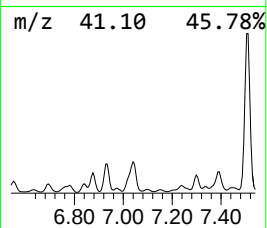
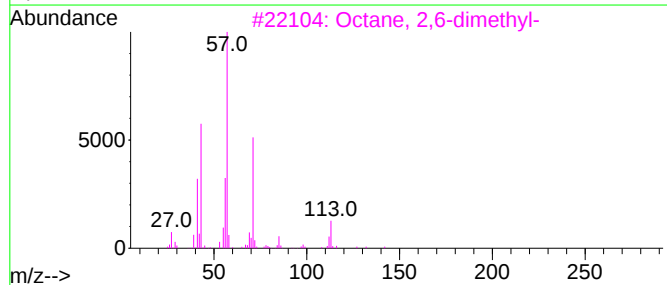
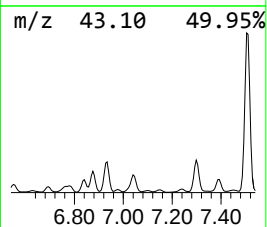
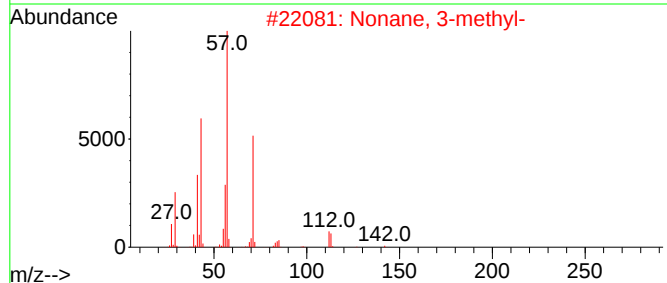
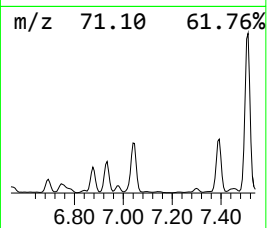
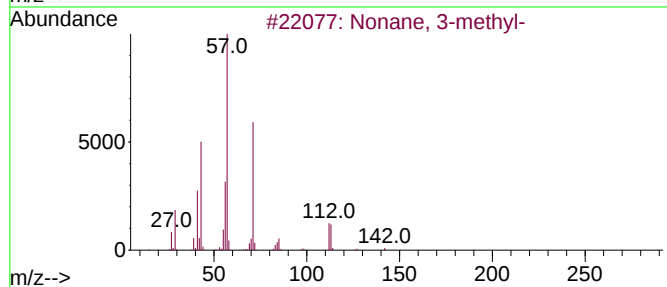
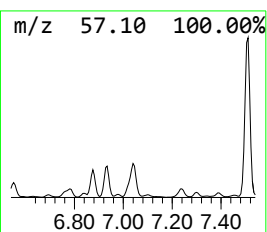
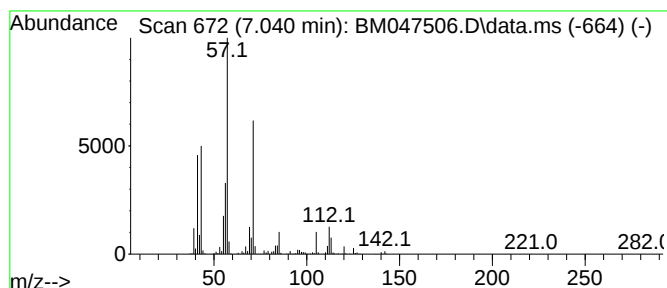
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.040	15.05 ng/ul	3838700	1,4-Dichlorobenzene-d4			7.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-		142	C10H22	005911-04-6	81
2	Nonane, 3-methyl-		142	C10H22	005911-04-6	76
3	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	64
4	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	64
5	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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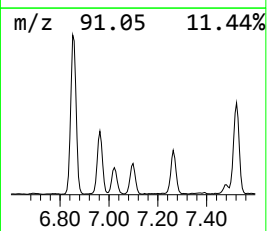
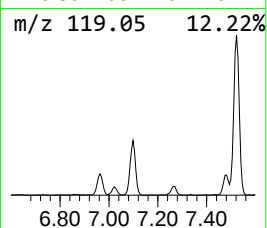
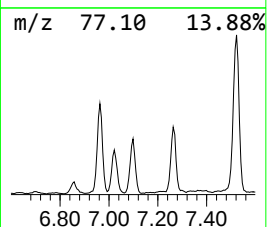
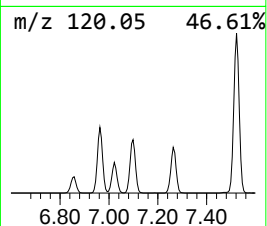
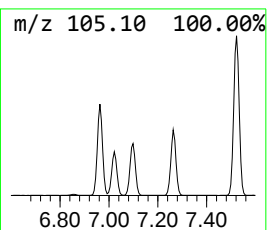
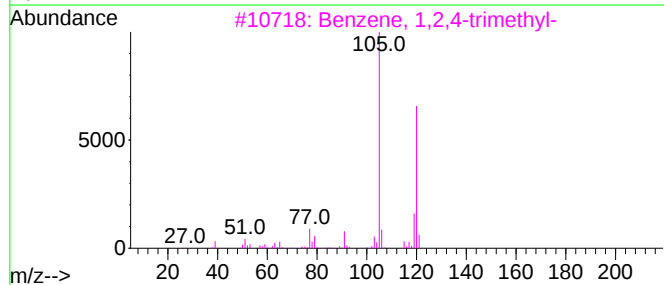
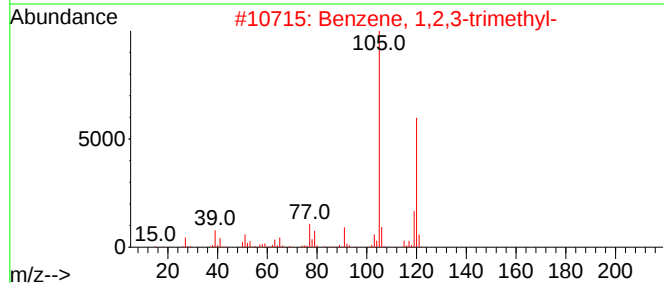
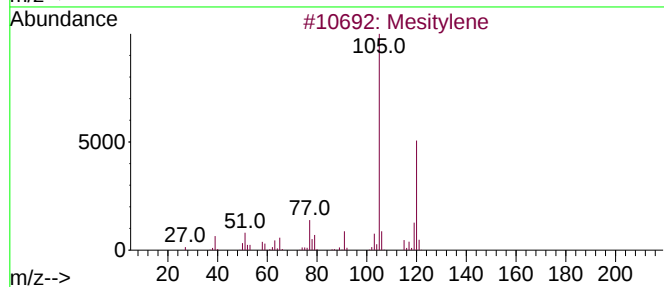
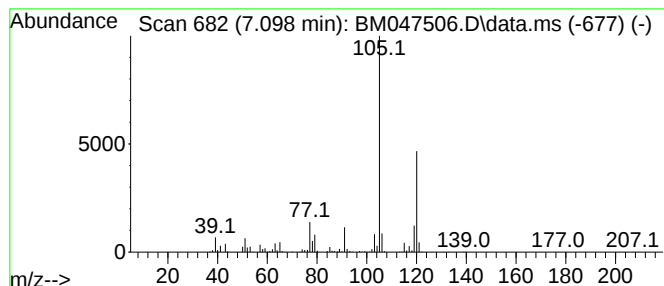
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Mesitylene Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.098	4.66 ng/ul	1188420	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Mesitylene	120	C9H12	000108-67-8	94
5			Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

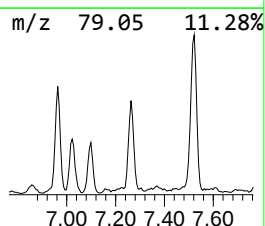
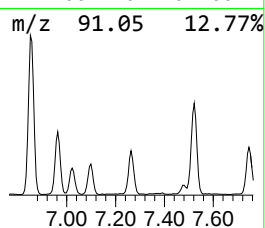
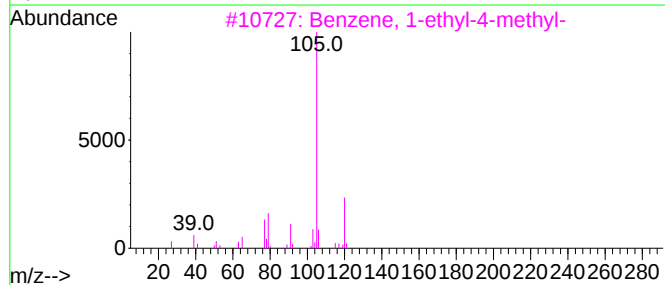
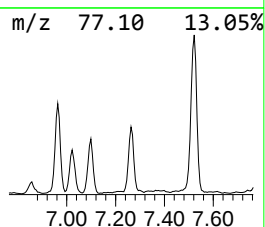
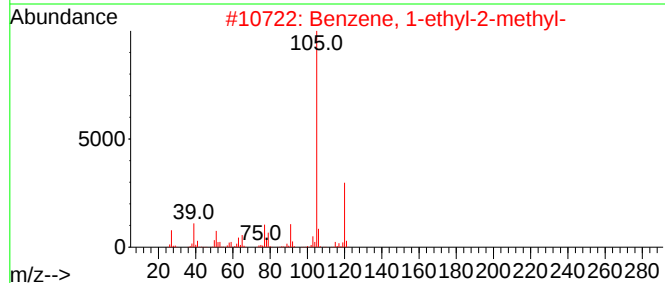
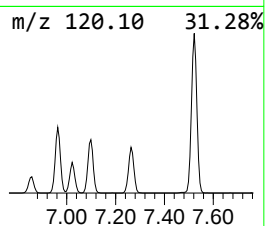
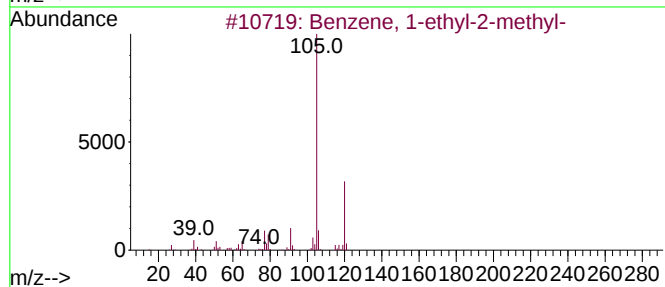
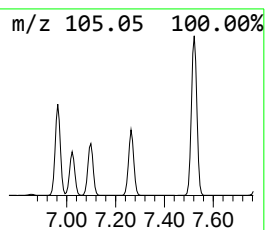
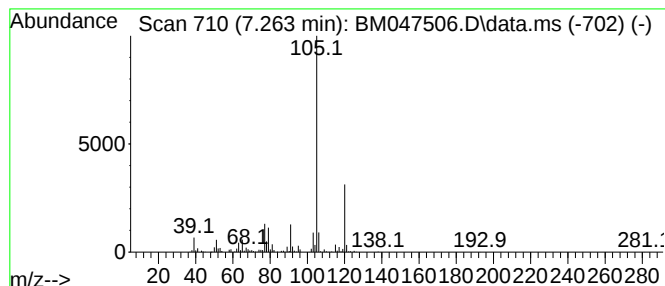
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzene, 1-ethyl-4-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.263	5.91 ng/ul	1508730	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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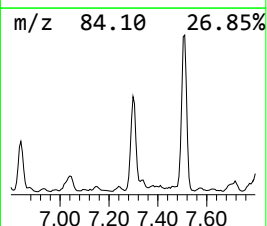
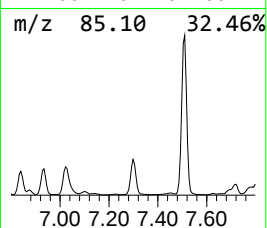
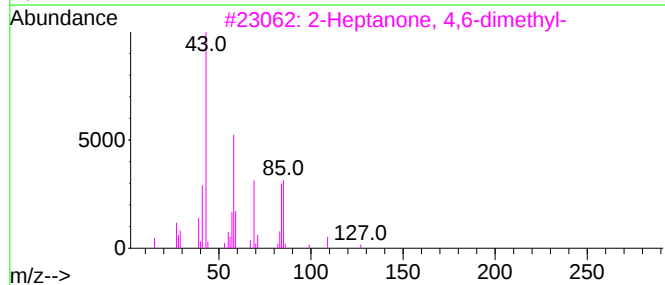
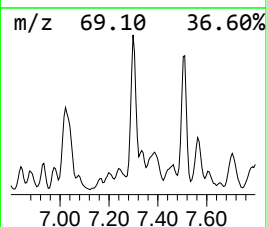
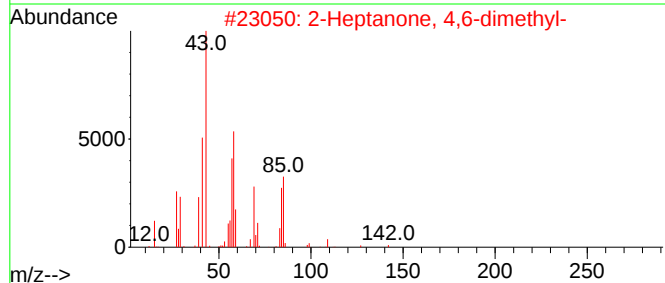
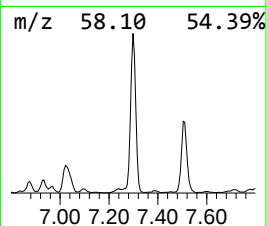
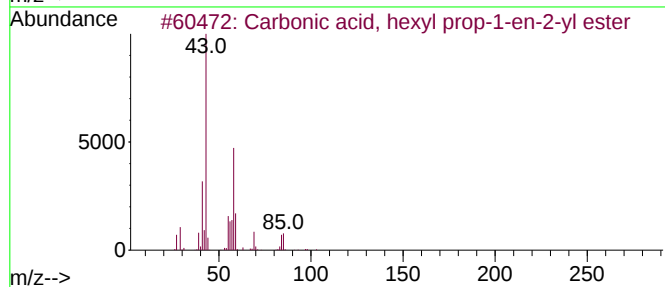
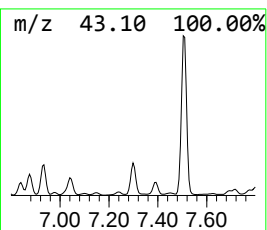
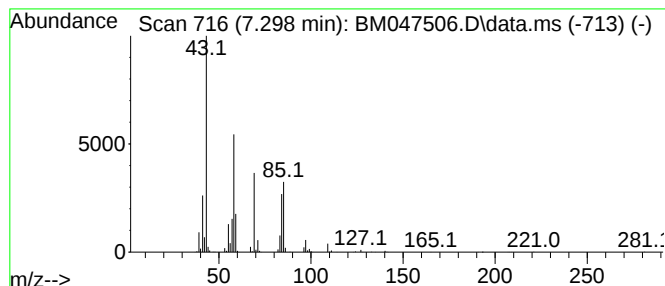
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Carbonic acid, hexyl prop-1... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.298	5.73 ng/ul	1460870	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, hexyl prop-1-en-2...	186	C10H18O3	1000382-54-2	59
2			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	56
3			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	45
4			2-Heptanone, 4-methyl-	128	C8H16O	006137-06-0	38
5			2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	38



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Sample : P3878-08DL2 20X
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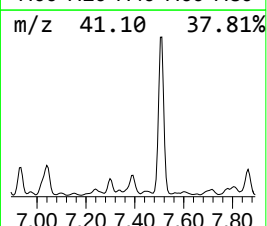
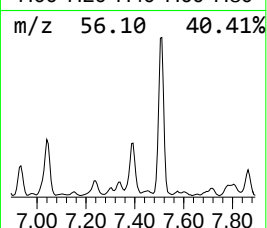
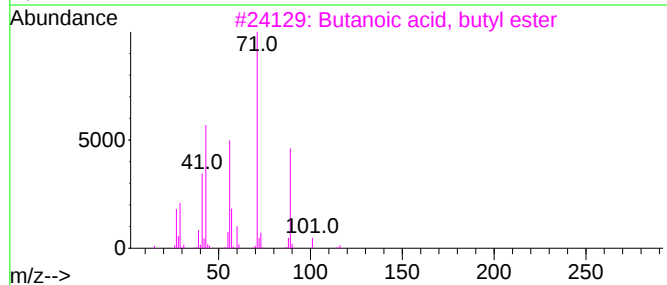
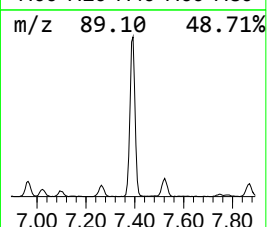
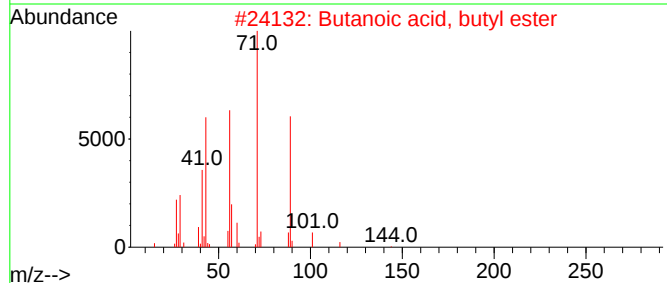
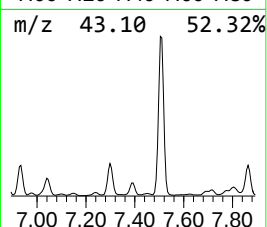
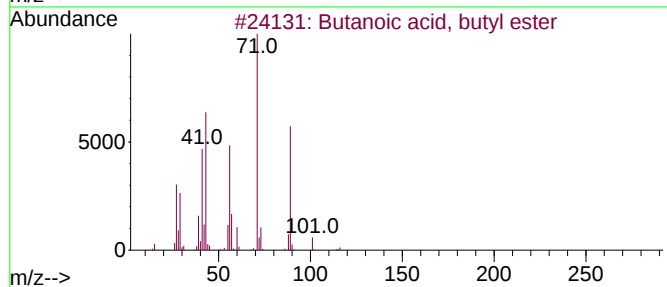
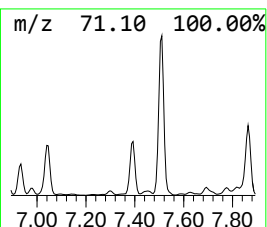
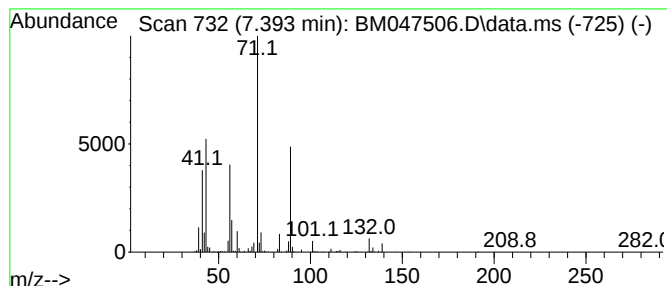
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Butanoic acid, butyl ester Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.393	7.86 ng/ul	2003730	1,4-Dichlorobenzene-d4	7.869	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	90
2	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	83
3	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	80
4	Butanoic acid, octyl ester	200	C12H24O2	000110-39-4	78
5	Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
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Instrument :
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ClientSampleId :
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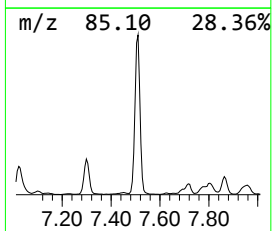
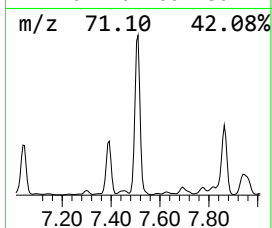
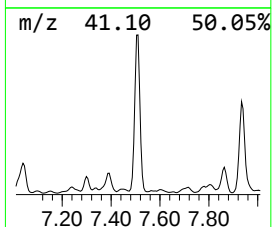
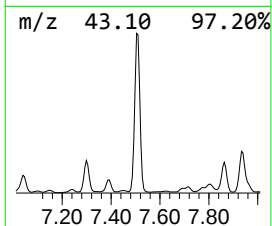
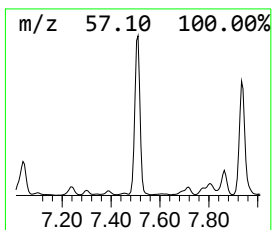
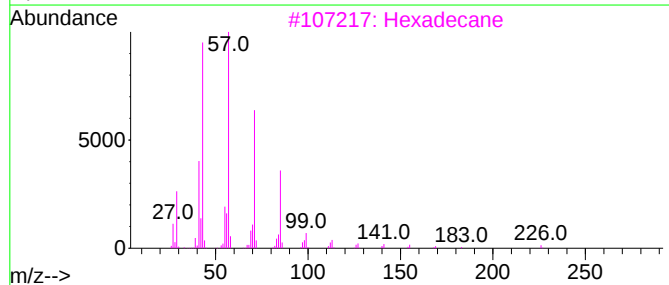
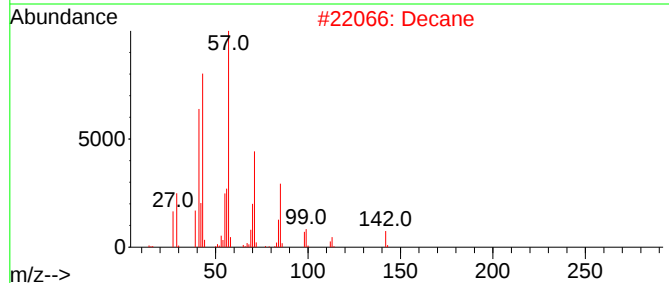
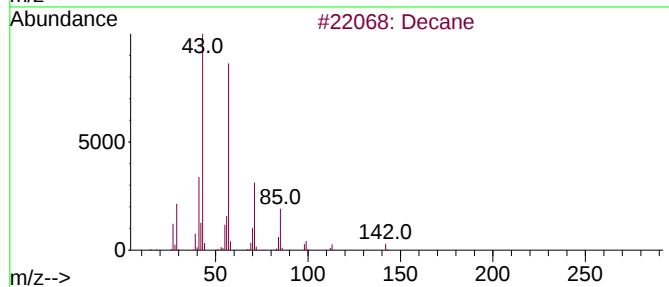
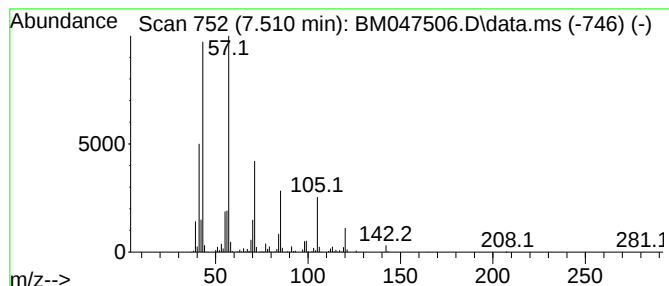
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.510	59.74 ng/ul	15238500	1,4-Dichlorobenzene-d4	7.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane	142	C10H22	000124-18-5	90
2		Decane	142	C10H22	000124-18-5	59
3		Hexadecane	226	C16H34	000544-76-3	53
4		Nonane	128	C9H20	000111-84-2	52
5		Tridecane	184	C13H28	000629-50-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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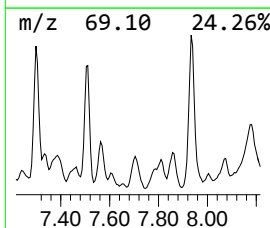
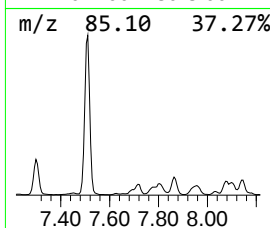
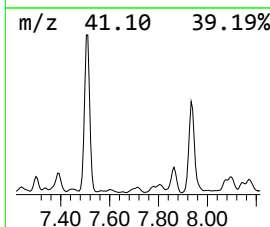
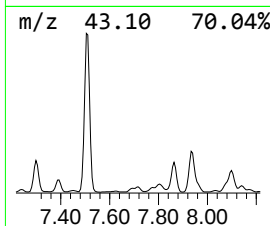
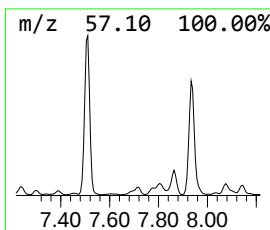
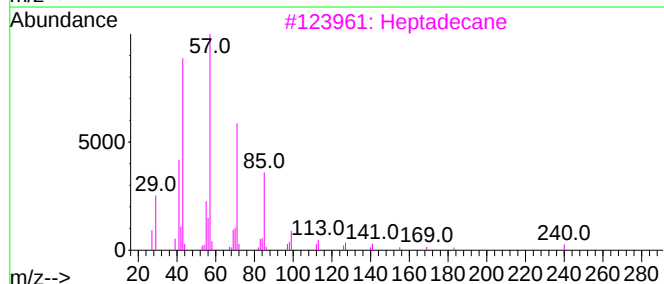
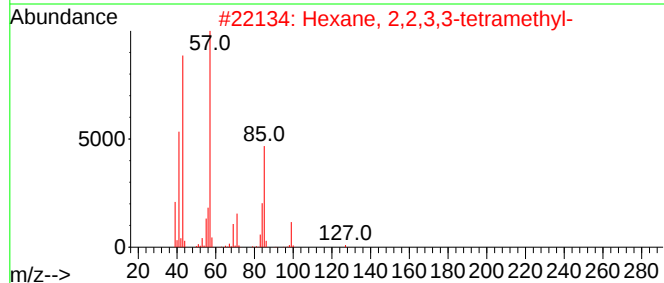
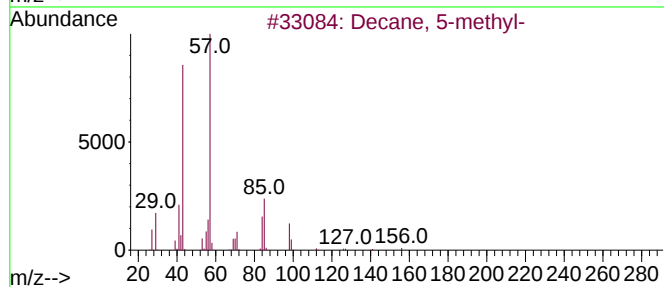
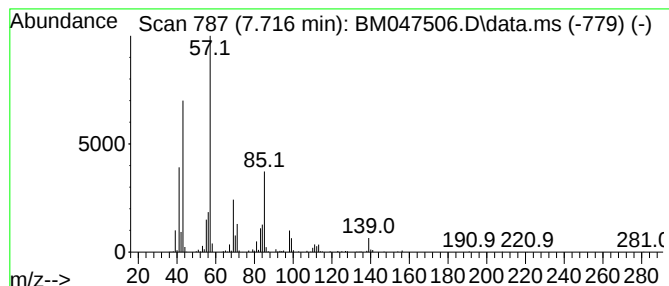
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.716	3.50 ng/ul	893548	1,4-Dichlorobenzene-d4	7.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5-methyl-	156	C11H24	013151-35-4	64
2		Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	43
3		Heptadecane	240	C17H36	000629-78-7	43
4		Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	43
5		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Sample : P3878-08DL2 20X
Misc :
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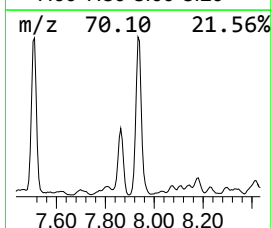
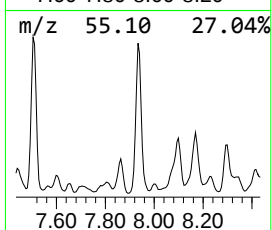
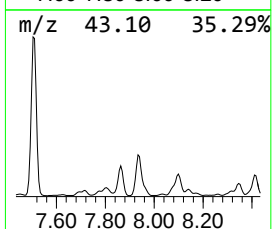
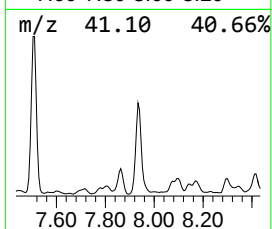
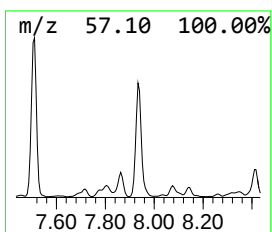
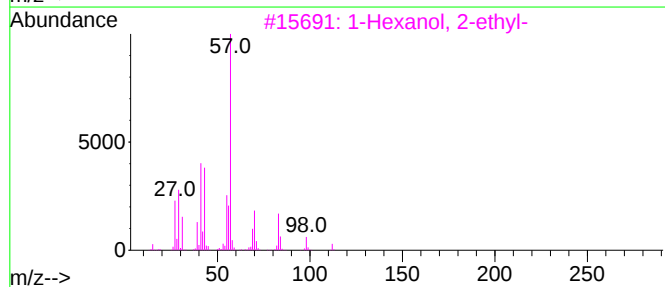
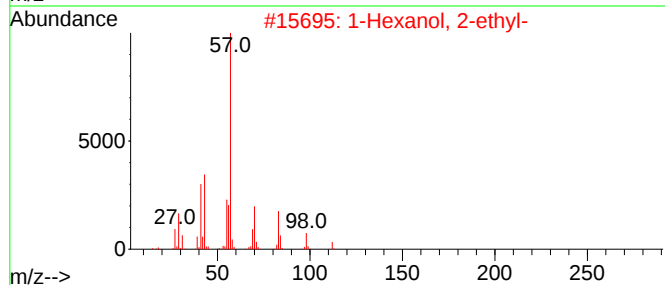
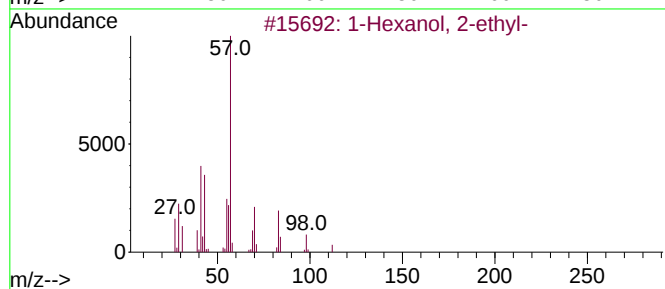
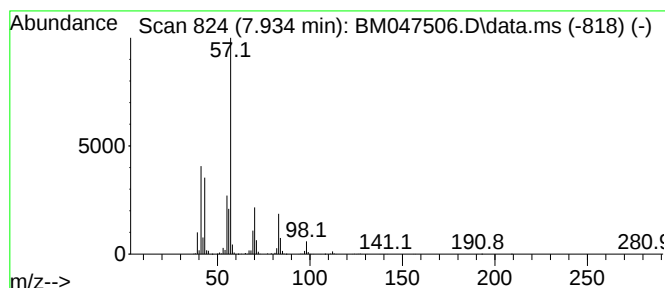
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.934	28.41 ng/ul	7247110	1,4-Dichlorobenzene-d4			7.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	74
5	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Sample : P3878-08DL2 20X
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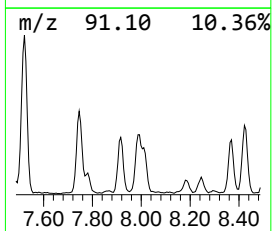
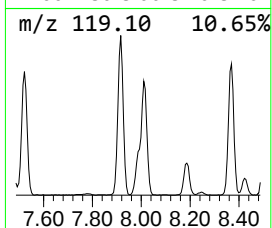
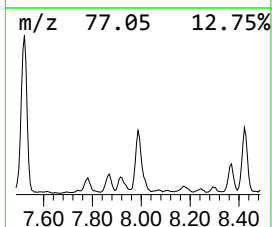
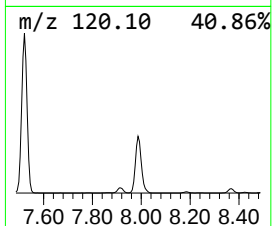
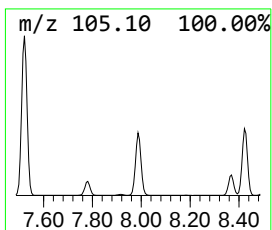
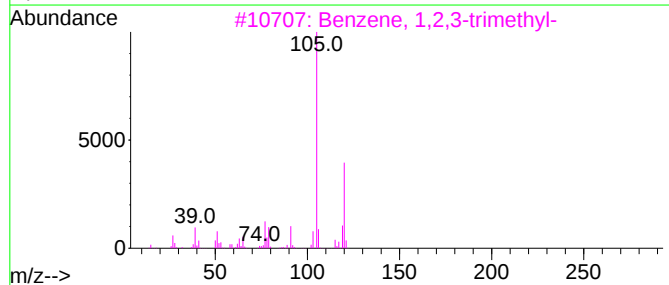
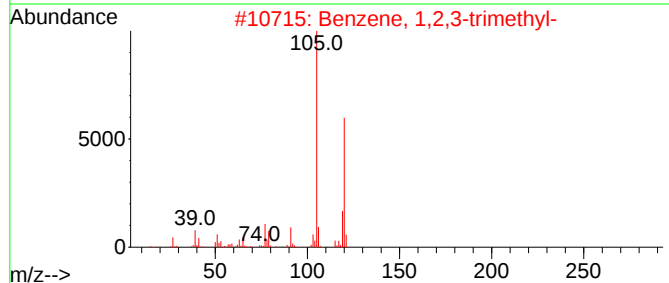
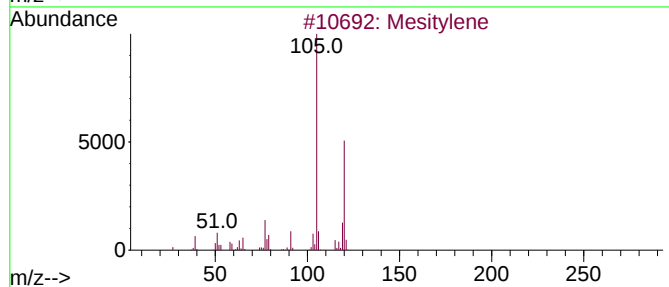
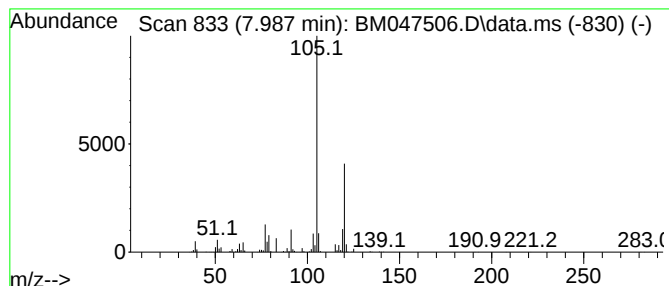
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Benzene, 1,2,3-trimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.987	5.66 ng/ul	1444280	1,4-Dichlorobenzene-d4	7.869	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
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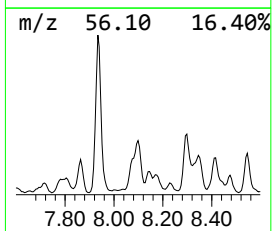
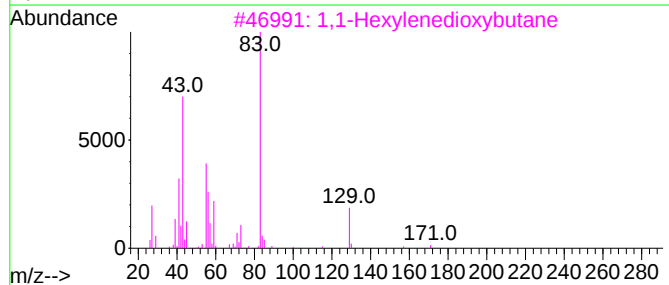
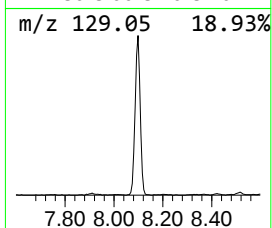
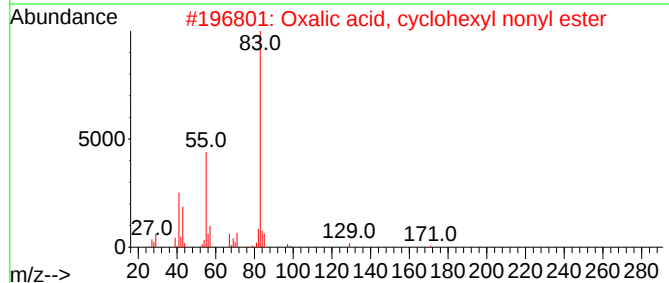
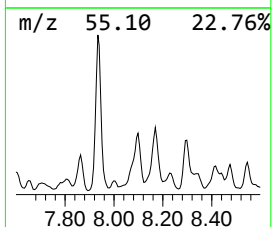
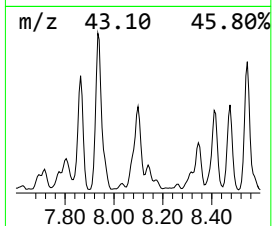
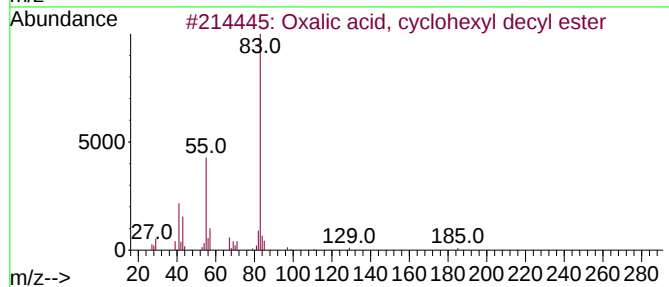
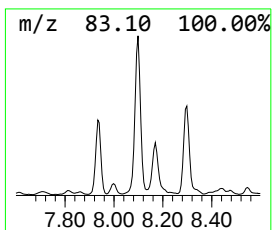
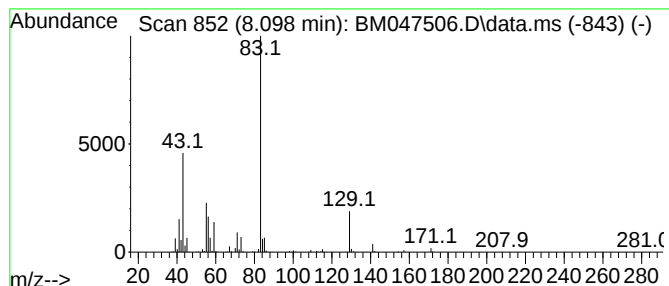
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Oxalic acid, cyclohexyl dec... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.098	12.70 ng/ul	3240390	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclohexyl decyl ester	312	C18H32O4	1000309-31-2	53
2			Oxalic acid, cyclohexyl nonyl ester	298	C17H30O4	1000309-31-1	53
3			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	43
4			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	42
5			Hexanal ethyl trans-2-hexenyl ac...	228	C14H28O2	1000431-42-6	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
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ALS Vial : 11 Sample Multiplier: 1

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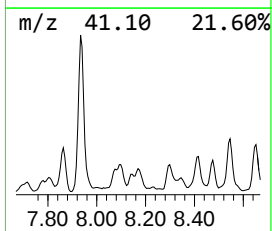
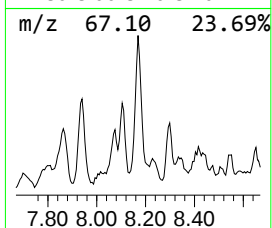
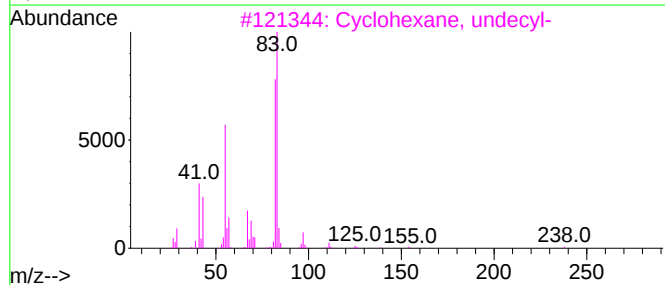
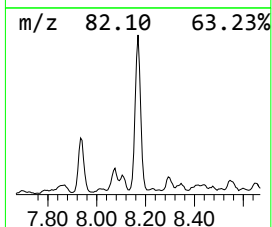
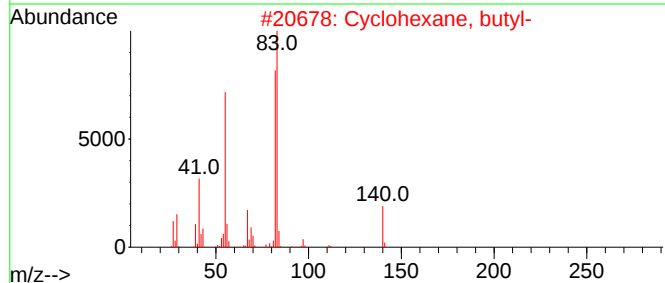
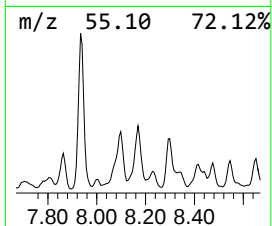
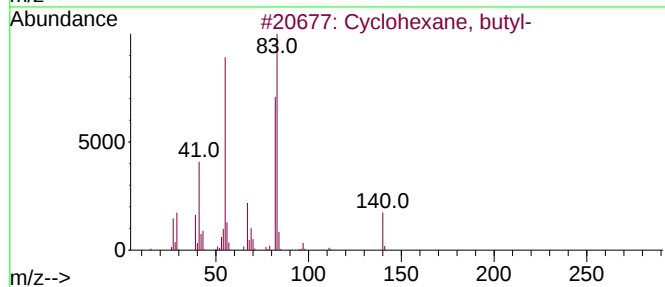
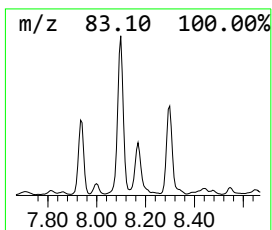
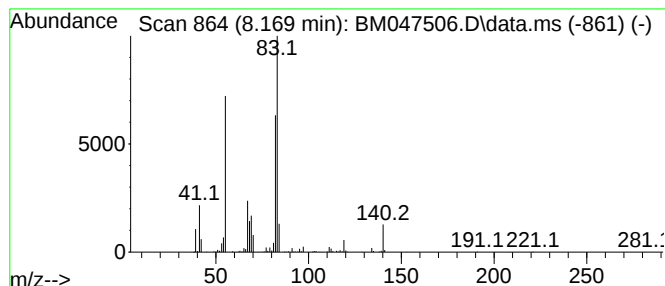
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 (DEL) Alkane: Cyclic8.169 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.169	5.00 ng/ul	1275220	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	86
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
3			Cyclohexane, undecyl-	238	C17H34	054105-66-7	72
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

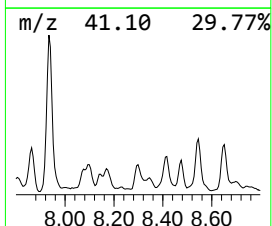
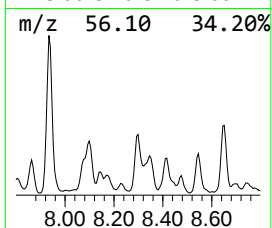
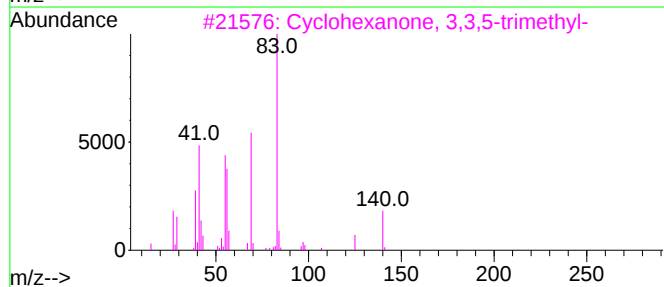
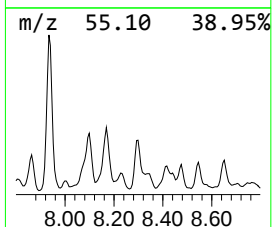
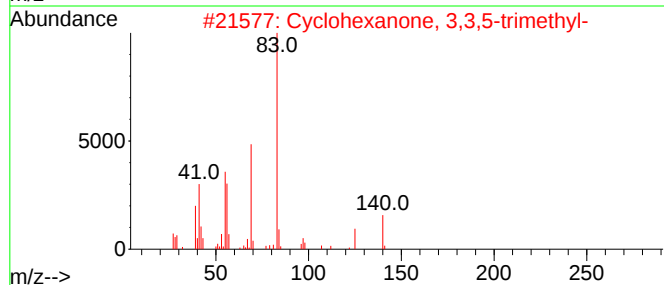
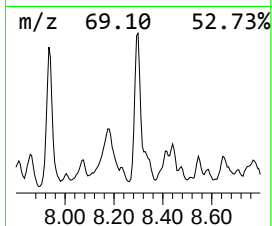
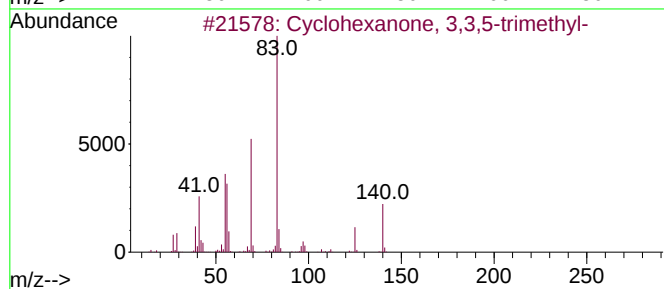
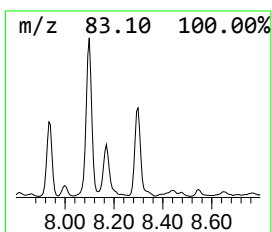
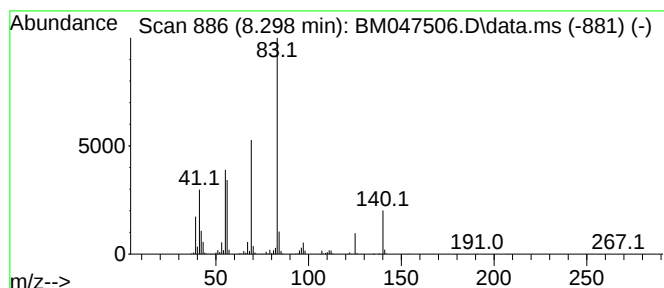
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 Cyclohexanone, 3,3,5-trimet... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.298	7.80 ng/ul	1988700	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5			Cyclopentanone, 3-butyl-	140	C9H16O	057283-81-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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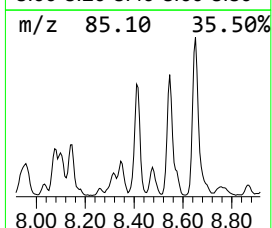
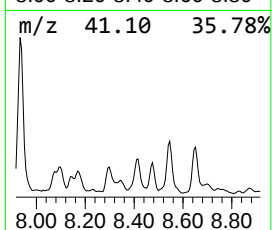
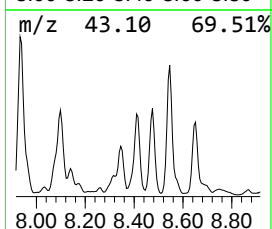
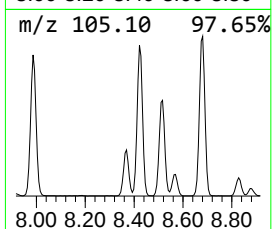
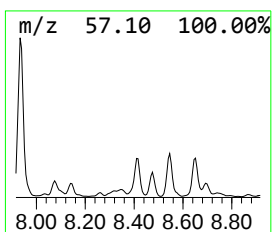
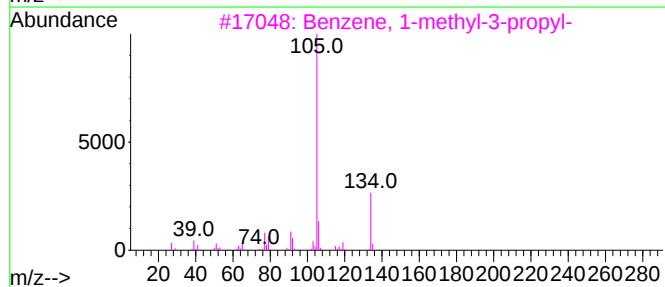
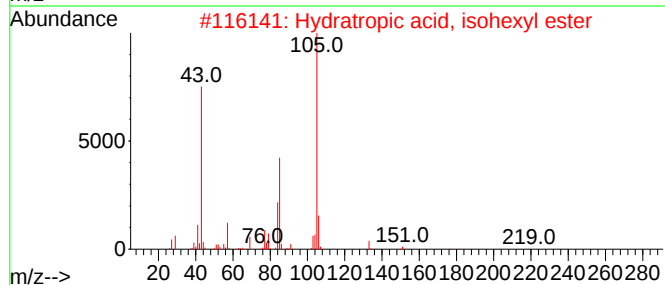
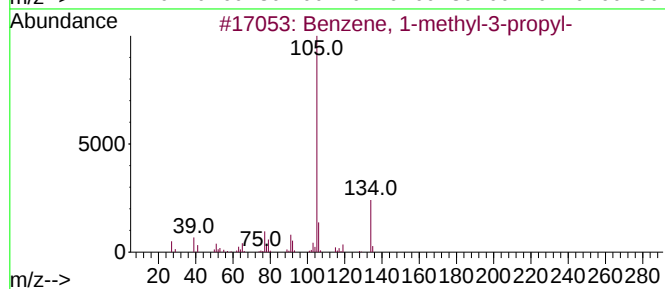
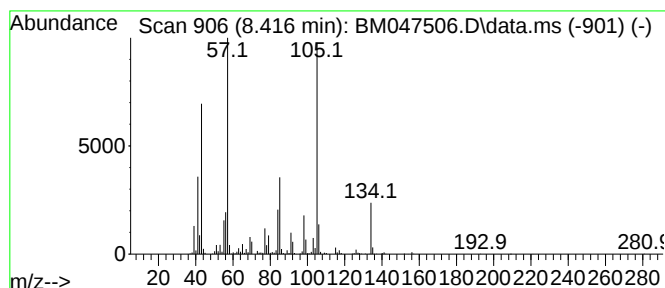
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 26 Benzene, 1-methyl-3-propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.416	10.76 ng/ul	2743580	1,4-Dichlorobenzene-d4	7.869	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	70
2	Hydratropic acid, isohexyl ester	234	C15H22O2	1000415-06-2	43
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	38
5	Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

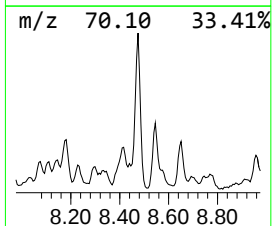
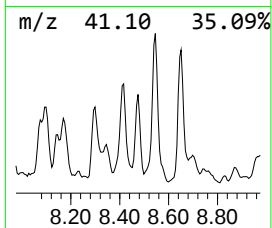
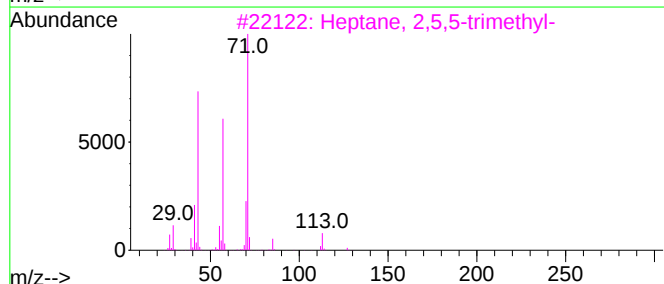
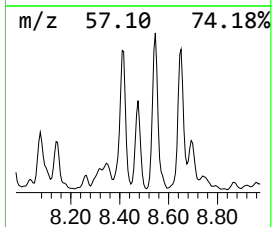
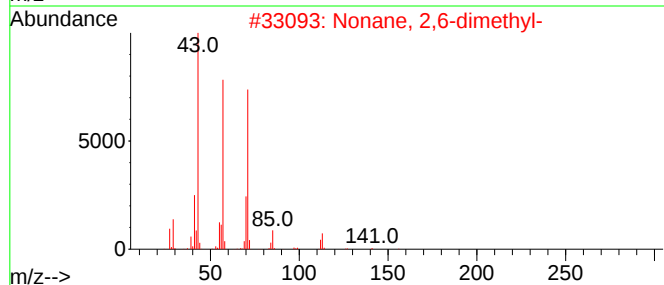
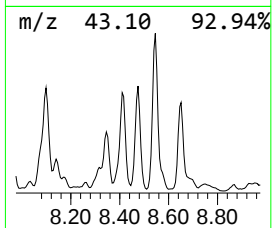
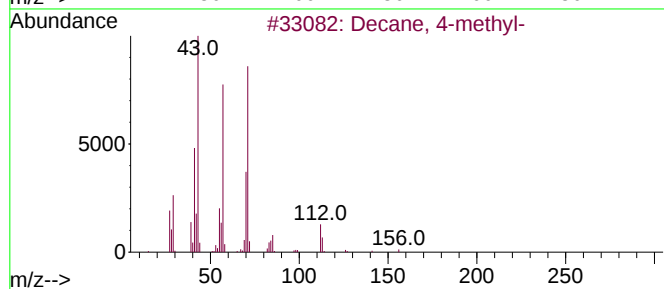
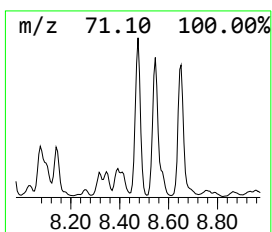
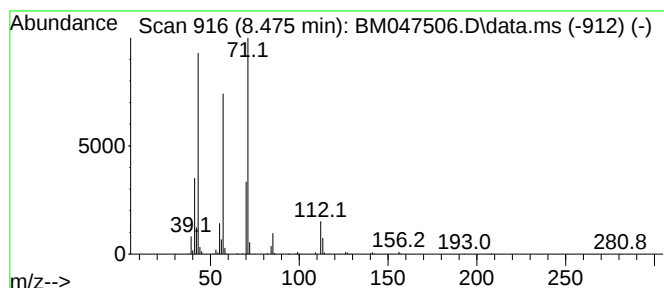
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.475	4.85 ng/ul	1236170	1,4-Dichlorobenzene-d4	7.869	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	90
2	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	83
3	Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	78
4	Decane, 4-methyl-	156	C11H24	002847-72-5	76
5	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Sample : P3878-08DL2 20X
Misc :
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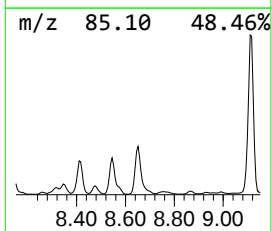
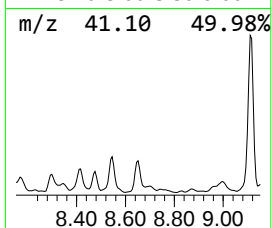
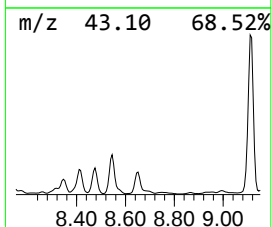
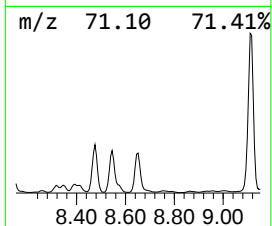
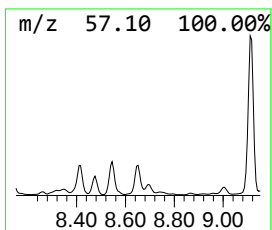
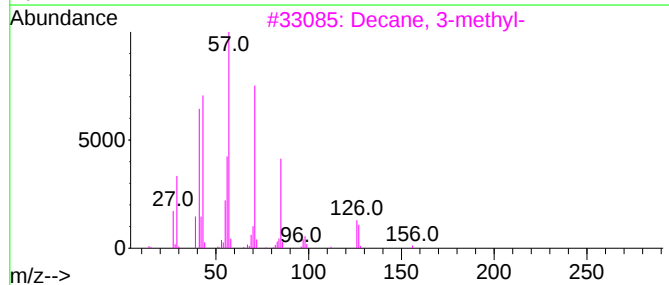
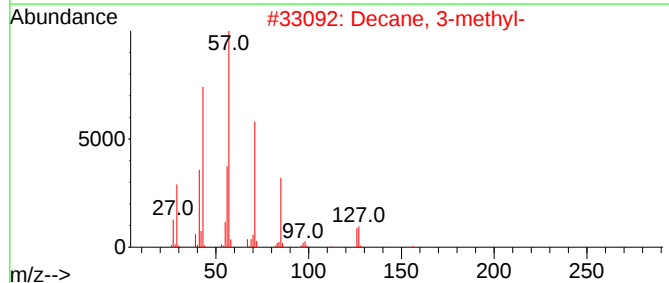
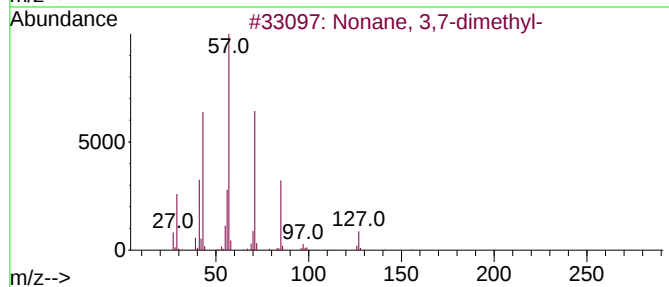
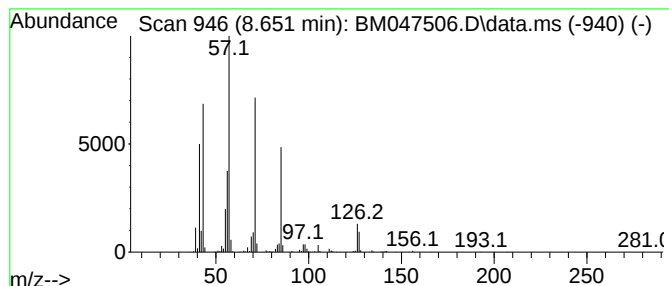
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.651	10.04 ng/ul	2560570	1,4-Dichlorobenzene-d4			7.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	87
2	Decane, 3-methyl-		156	C11H24	013151-34-3	87
3	Decane, 3-methyl-		156	C11H24	013151-34-3	83
4	Nonane, 5-(2-methylpropyl)-		184	C13H28	062185-53-9	80
5	Dodecane, 3-methyl-		184	C13H28	017312-57-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

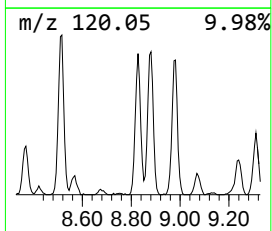
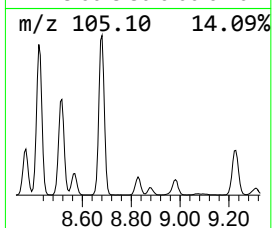
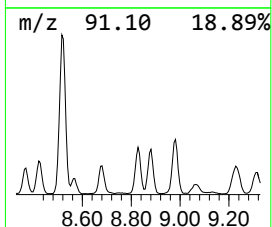
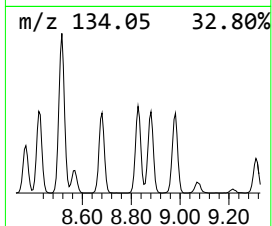
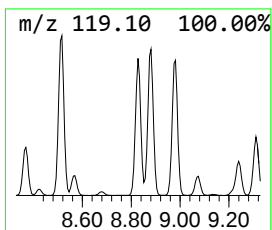
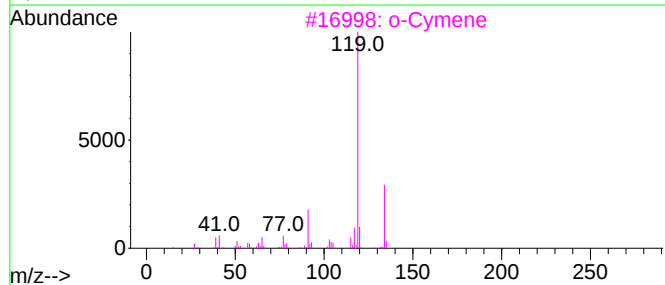
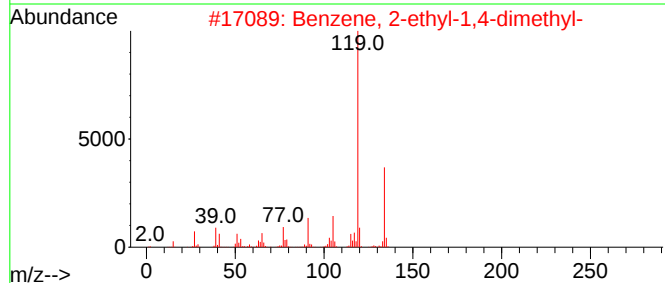
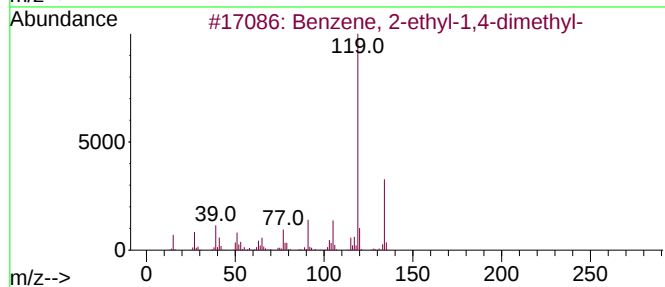
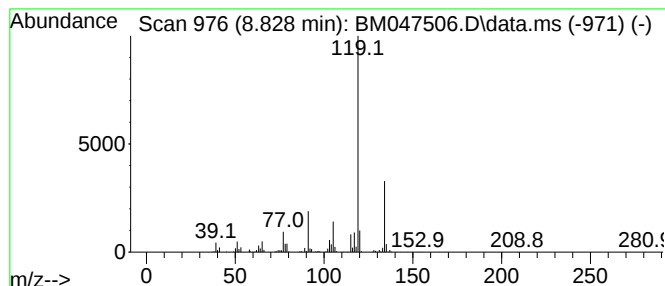
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.828	4.24 ng/ul	1080270	1,4-Dichlorobenzene-d4			7.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-		134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-		134	C10H14	001758-88-9	95
3	o-Cymene		134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	95
5	Benzene, 2-ethyl-1,4-dimethyl-		134	C10H14	001758-88-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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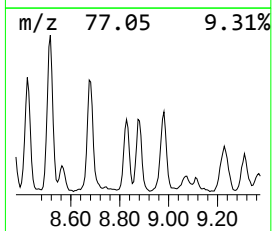
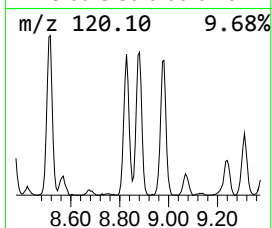
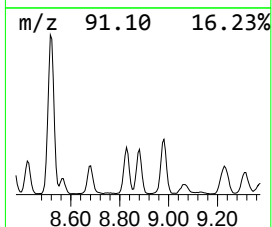
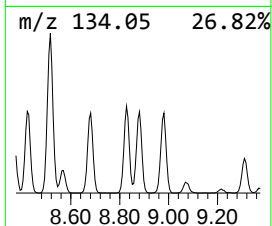
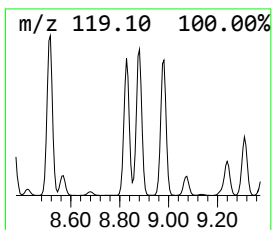
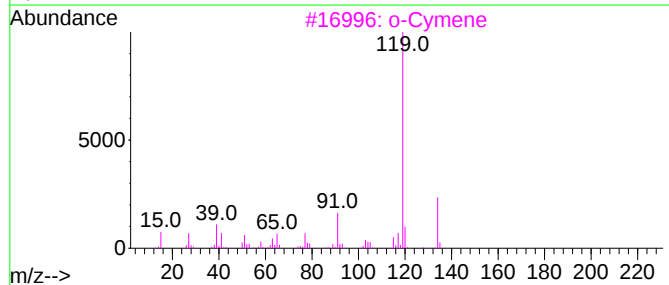
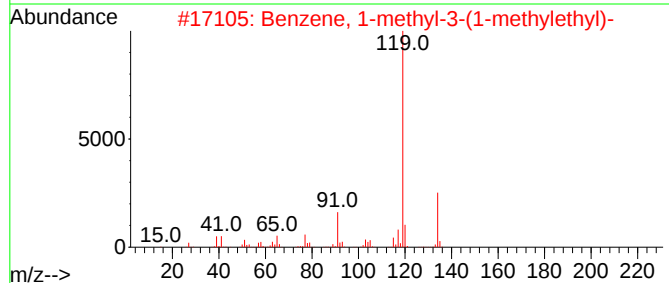
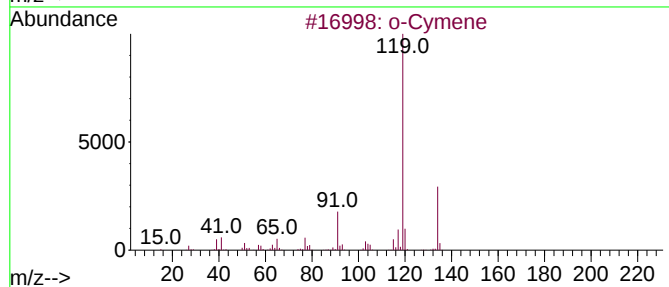
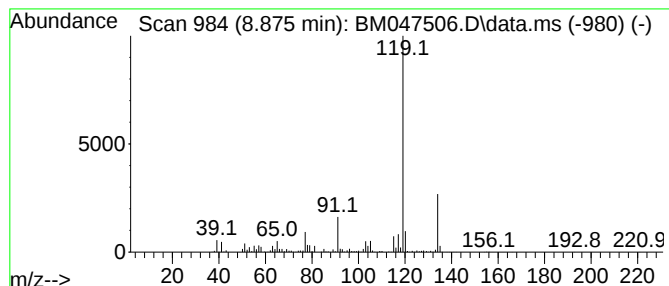
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 o-Cymene Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.875	5.31 ng/ul	1355430	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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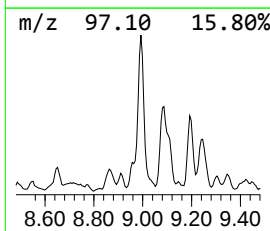
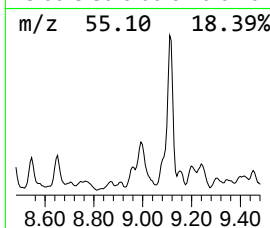
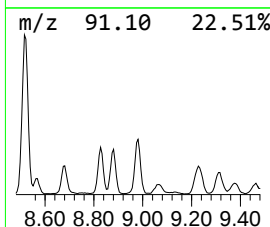
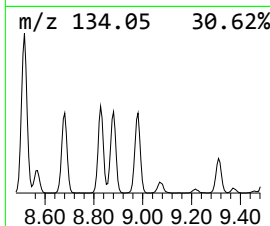
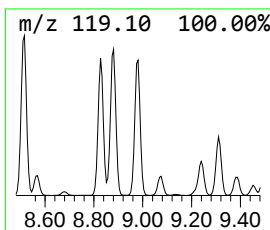
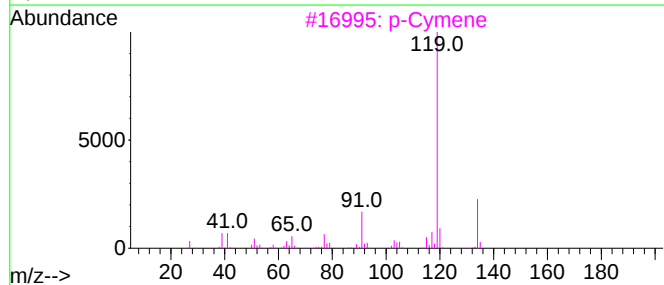
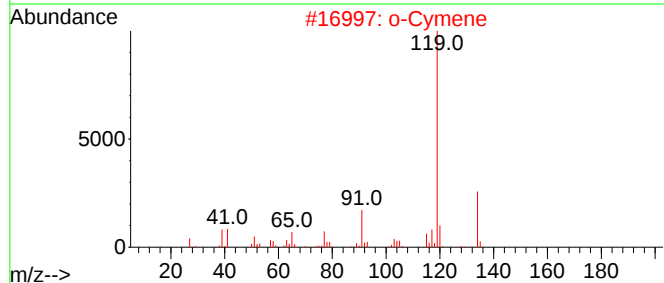
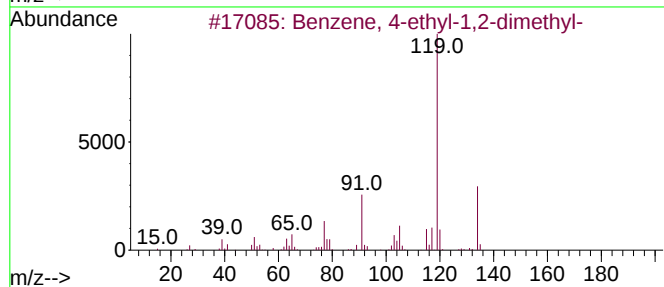
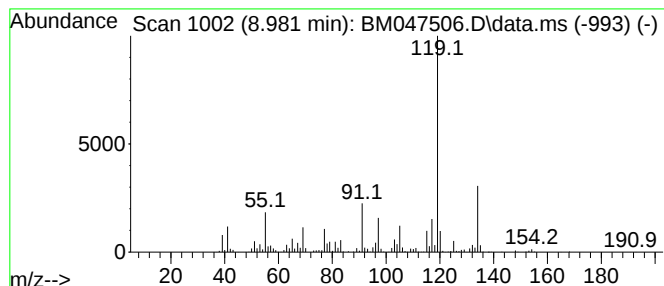
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.981	12.32 ng/ul	3141800	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

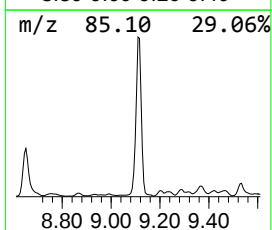
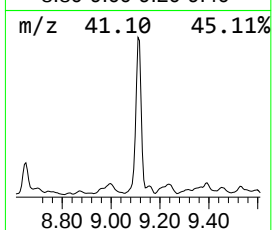
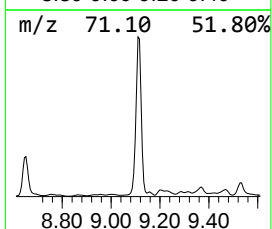
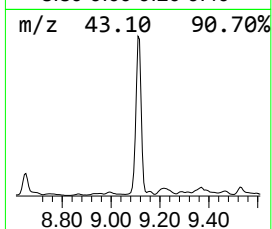
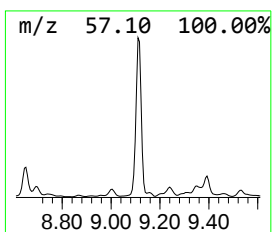
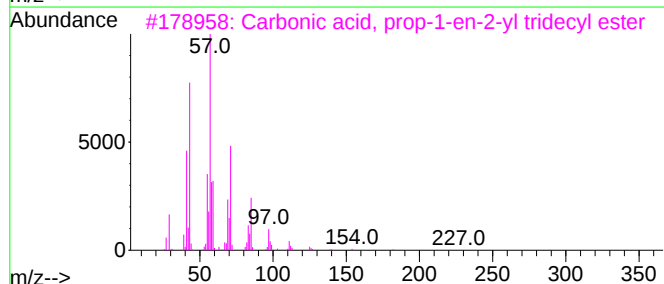
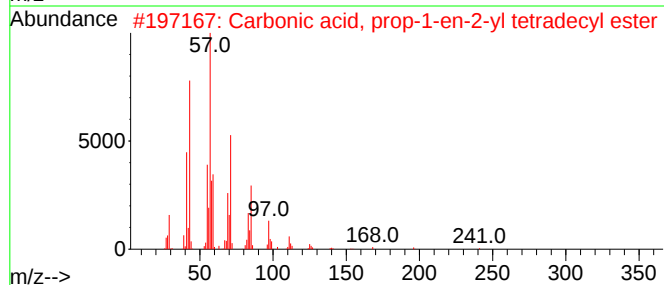
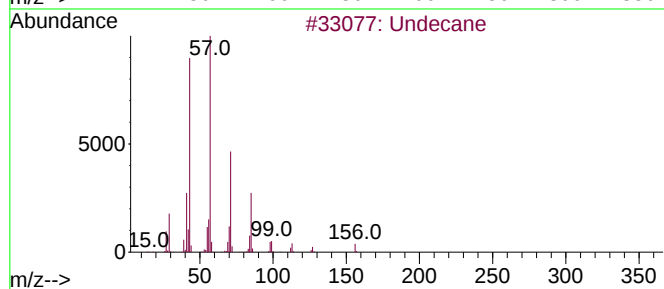
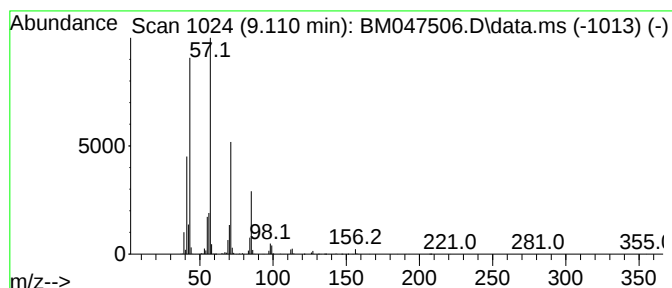
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.110	44.23 ng/ul	11281200	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	90
4			Hexadecane	226	C16H34	000544-76-3	86
5			Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

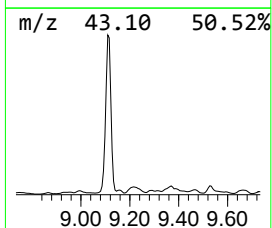
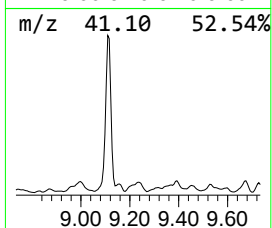
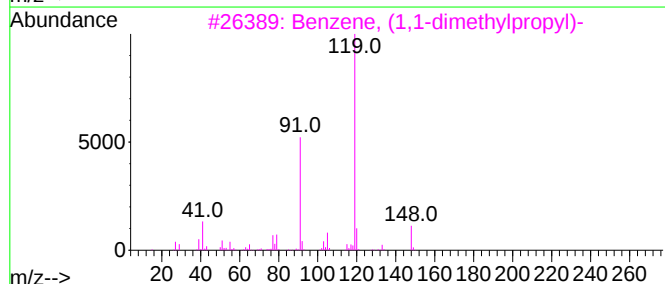
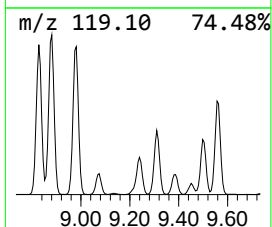
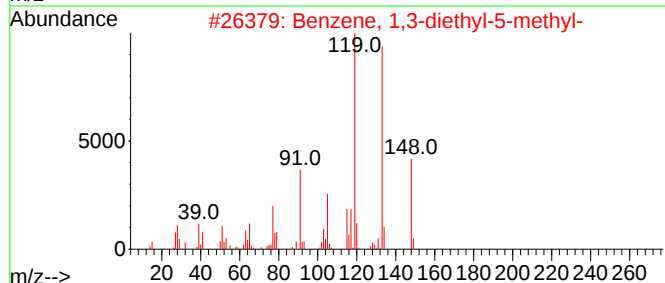
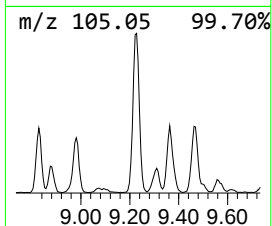
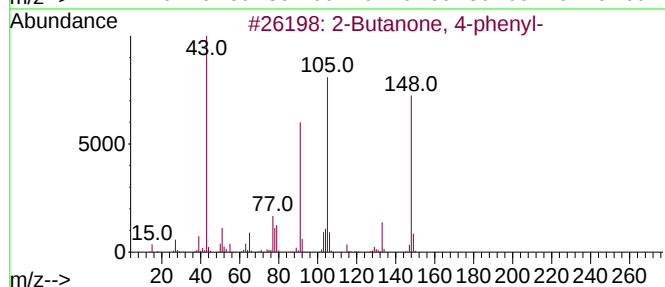
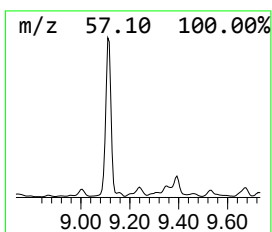
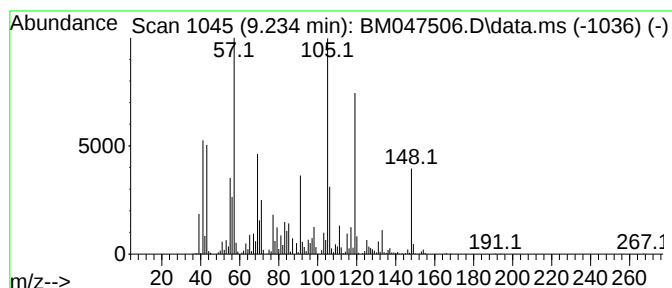
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 2-Butanone, 4-phenyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.234	9.58 ng/ul	2444610	1,4-Dichlorobenzene-d4	7.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	70	
2	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	64	
3	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	47	
4	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	42	
5	Benzene, 4-(chloromethyl)-1,2-di...	154	C9H11Cl	000102-46-5	41	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

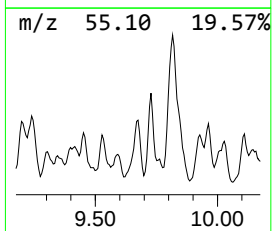
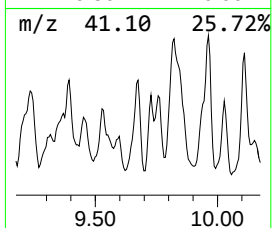
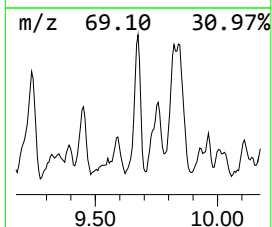
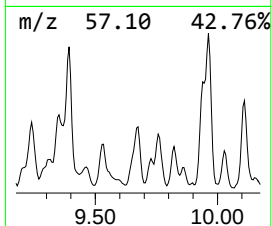
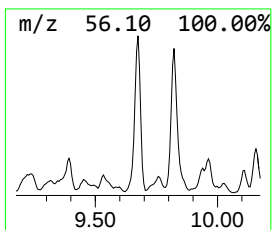
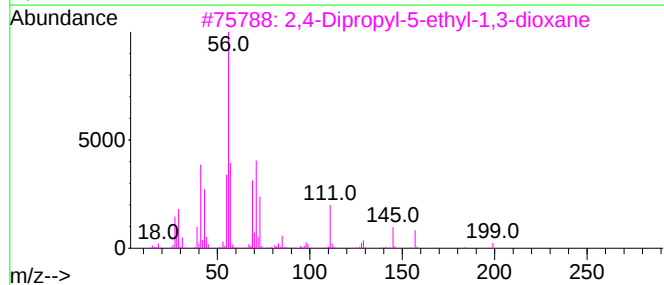
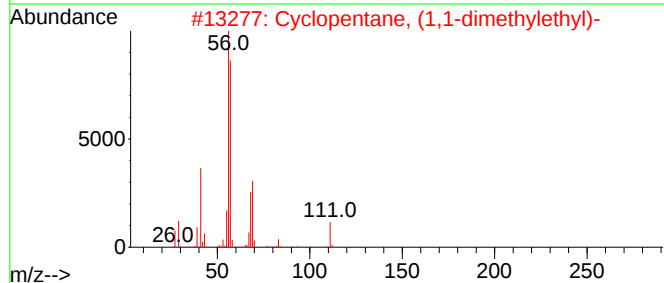
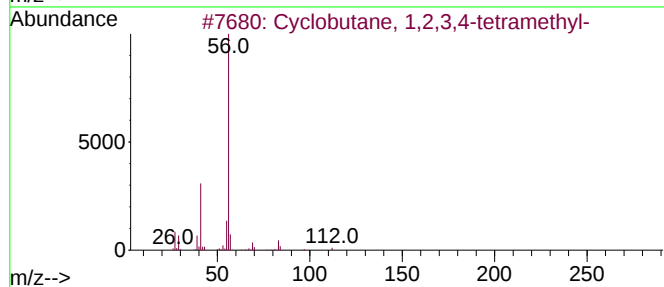
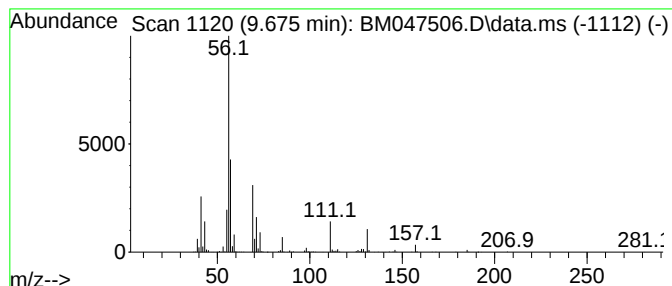
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 (DEL) Alkane: Cyclic9.675 Concentration Rank 78

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.675	2.05 ng/ul	1340540	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	43
2		Cyclopentane, (1,1-dimethylethyl)-	126	C9H18	003875-52-3	38
3		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	16
4		2,4-Dimethyl-1-hexene	112	C8H16	016746-87-5	10
5		2-Propen-1-amine	57	C3H7N	000107-11-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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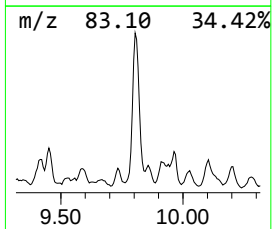
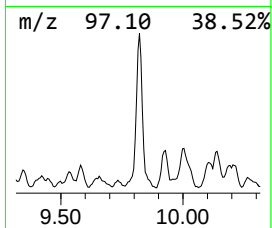
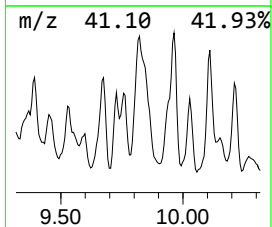
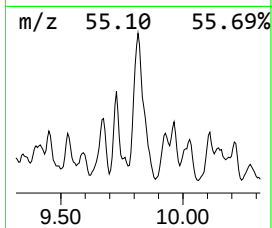
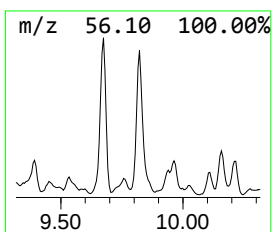
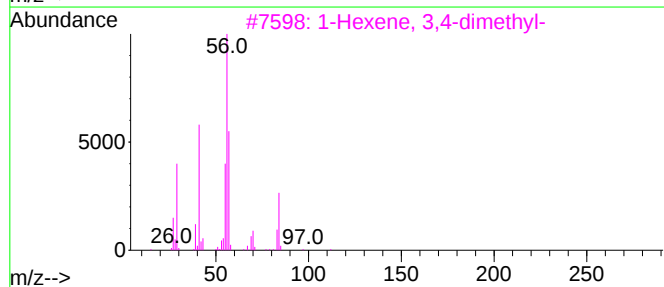
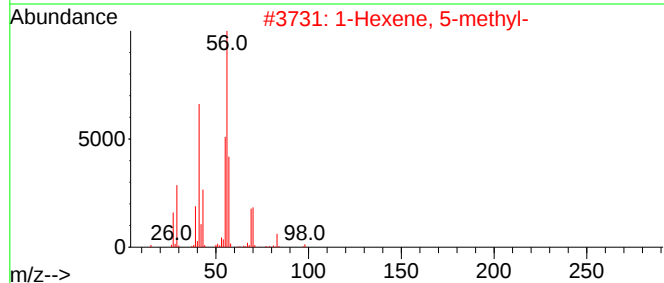
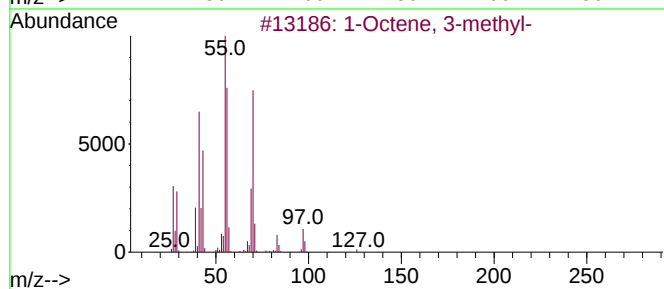
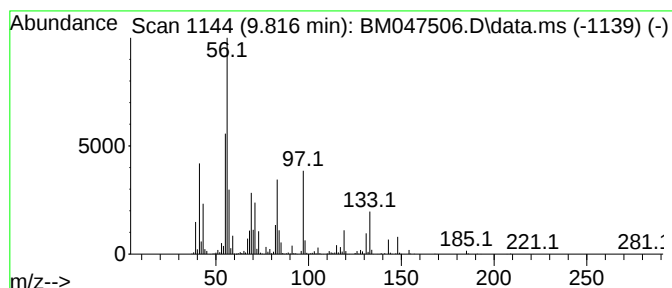
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 70

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.816	2.94 ng/ul	1924970	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Octene, 3-methyl-	126	C9H18	013151-08-1	35
2		1-Hexene, 5-methyl-	98	C7H14	003524-73-0	27
3		1-Hexene, 3,4-dimethyl-	112	C8H16	016745-94-1	27
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	27
5		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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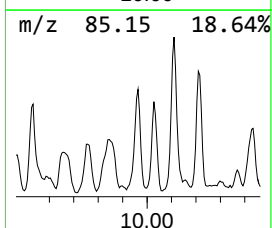
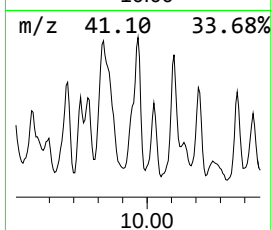
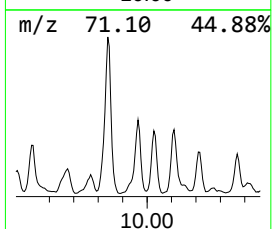
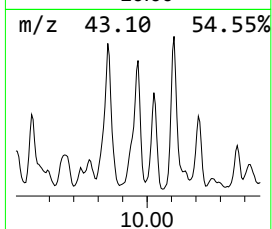
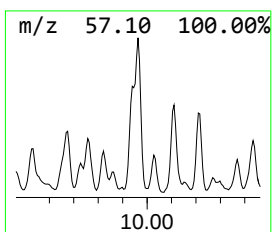
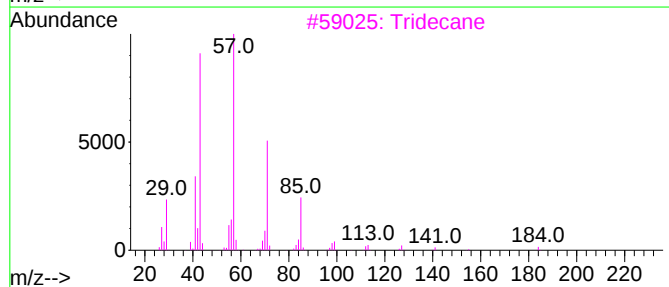
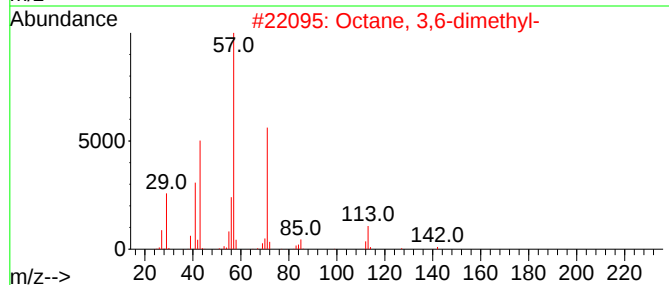
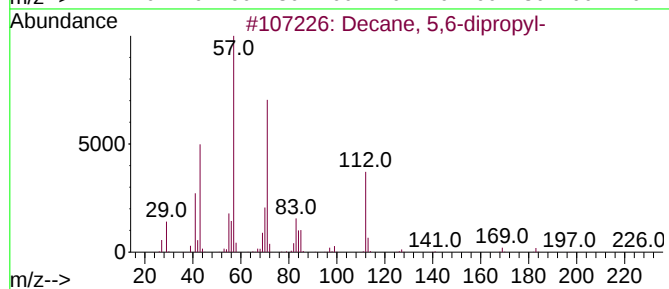
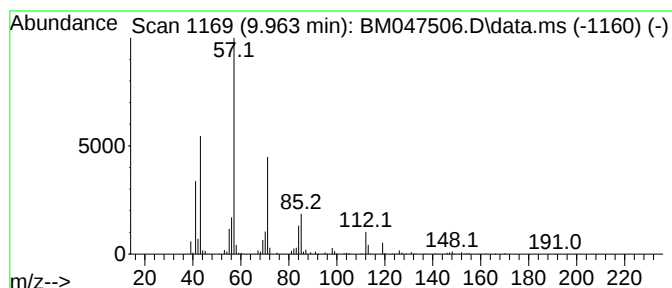
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	3.11 ng/ul	2037370	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	64
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	53
3		Tridecane	184	C13H28	000629-50-5	53
4		Octane, 4-ethyl-	142	C10H22	015869-86-0	53
5		1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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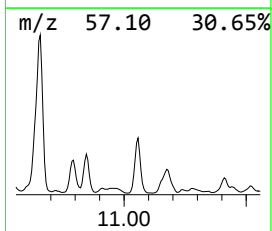
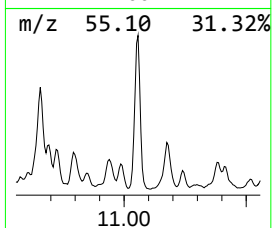
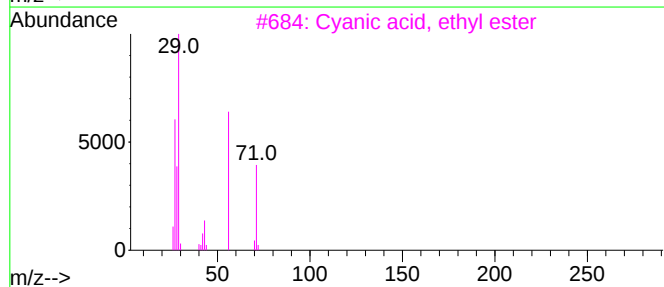
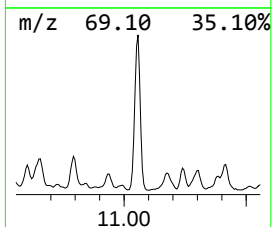
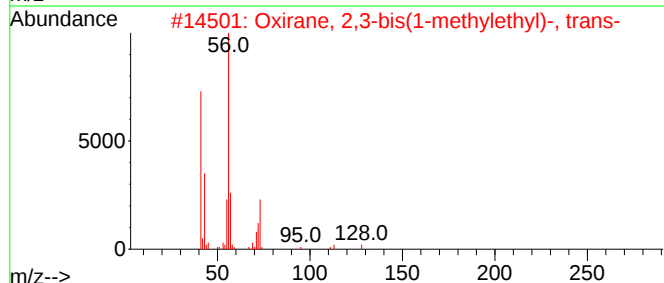
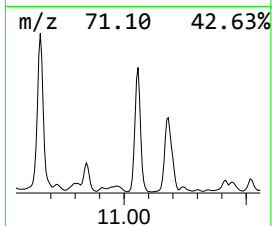
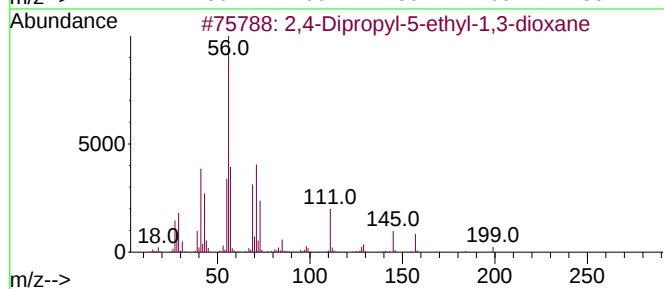
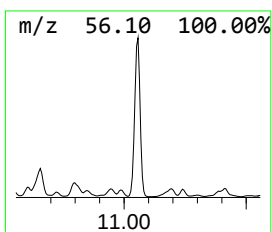
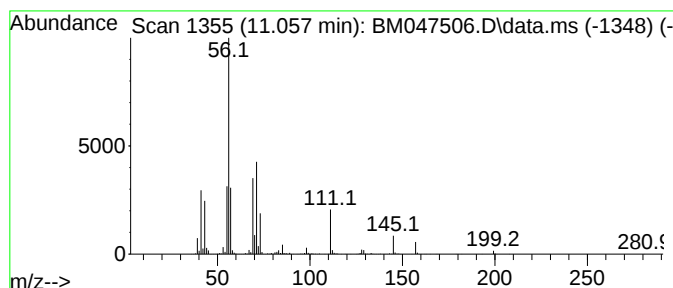
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.057	8.88 ng/ul	5811820	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	64
2		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	43
3		Cyanoic acid, ethyl ester	71	C3H5NO	000627-48-5	42
4		Cyclohexylamine	99	C6H13N	000108-91-8	35
5		Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

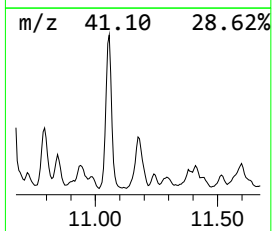
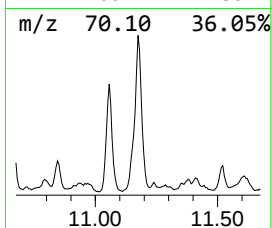
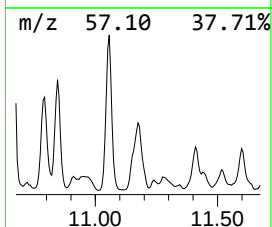
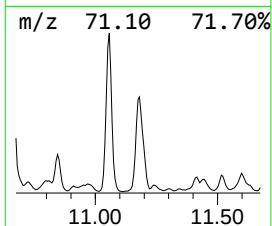
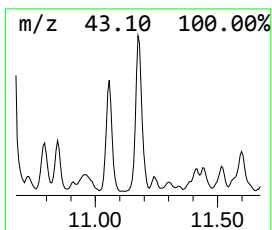
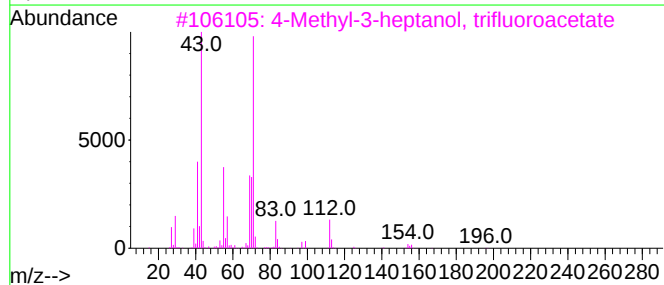
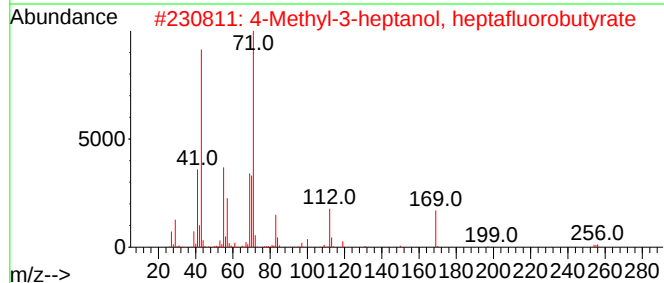
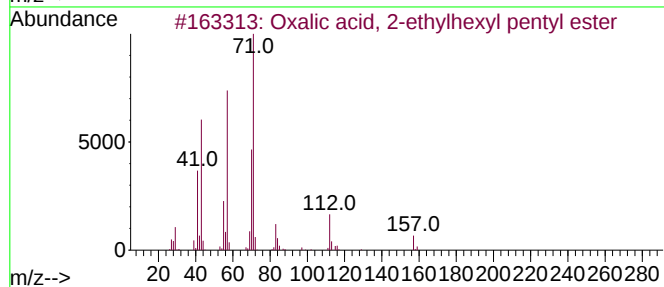
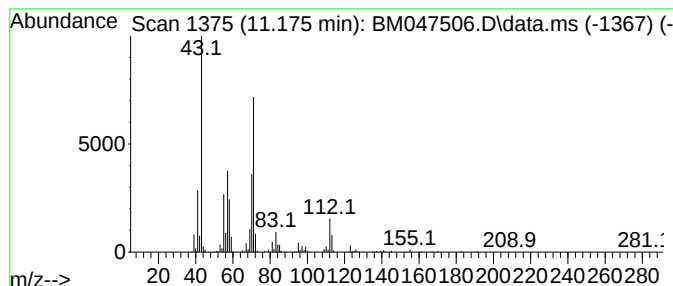
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 Oxalic acid, 2-ethylhexyl p... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.175	4.22 ng/ul	2762490	Naphthalene-d8	10.663	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	59
2	4-Methyl-3-heptanol, heptafluoro...	326	C12H17F7O2	1000365-51-5	53
3	4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	53
4	Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	47
5	2,3'-Bifuran, octahydro-	142	C8H14O2	073373-15-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
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Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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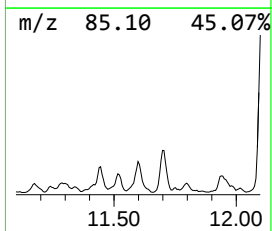
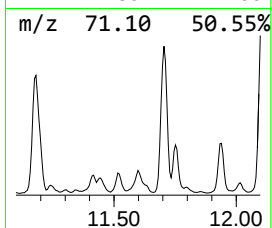
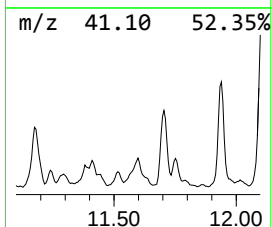
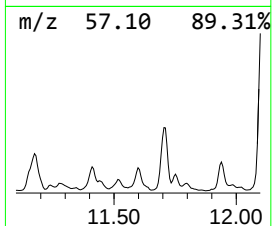
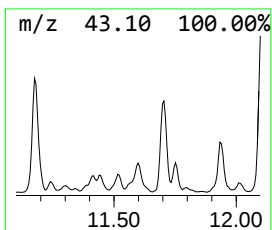
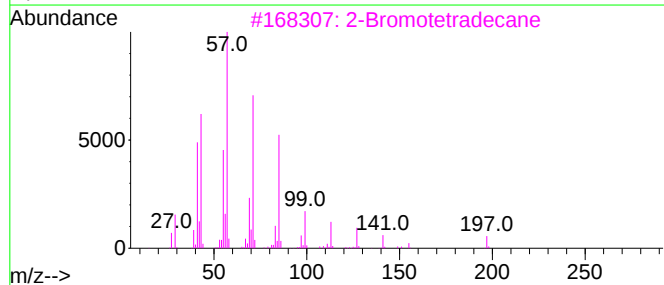
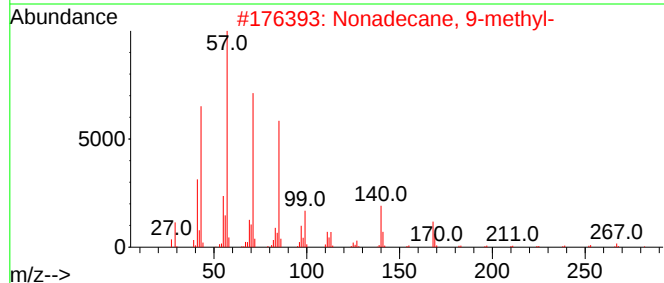
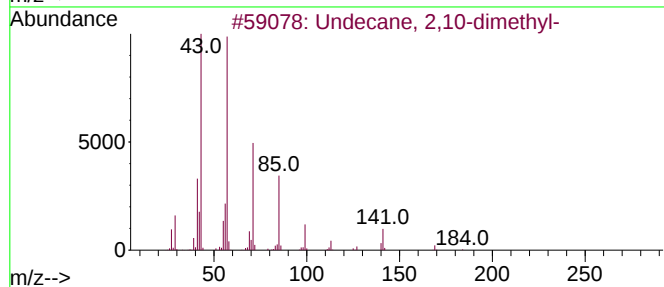
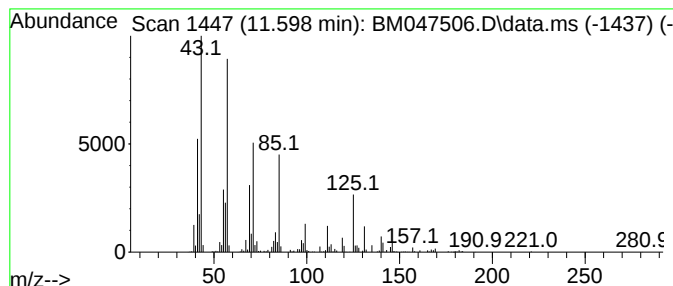
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 75

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.598	2.52 ng/ul	1648830	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 2,10-dimethyl-	184	C13H28	017301-27-8	70
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	58
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	58
4		Tridecane, 7-propyl-	226	C16H34	055045-09-5	58
5		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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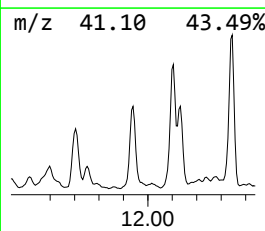
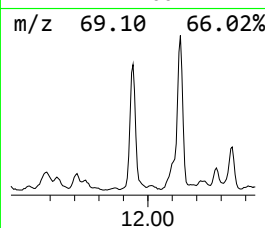
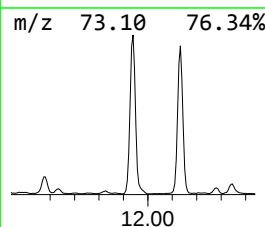
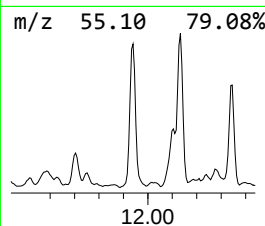
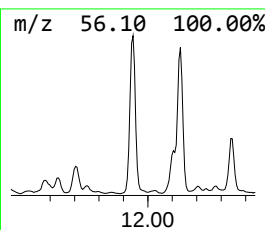
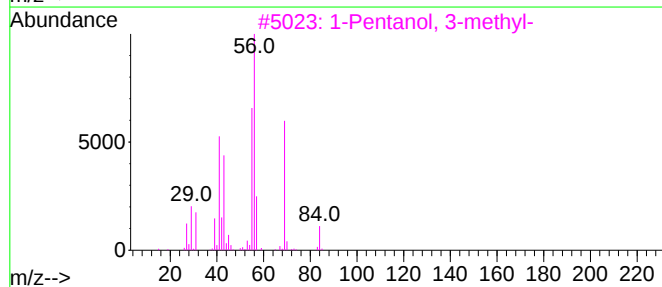
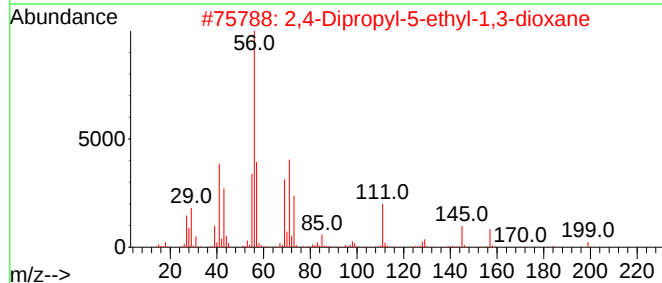
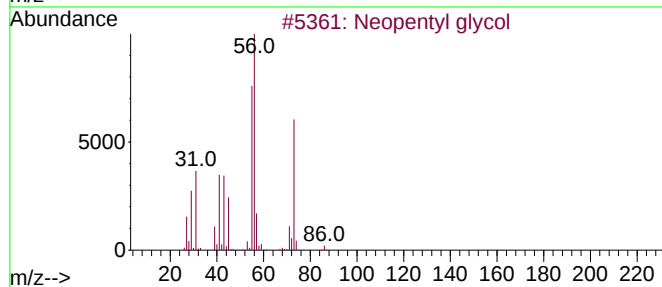
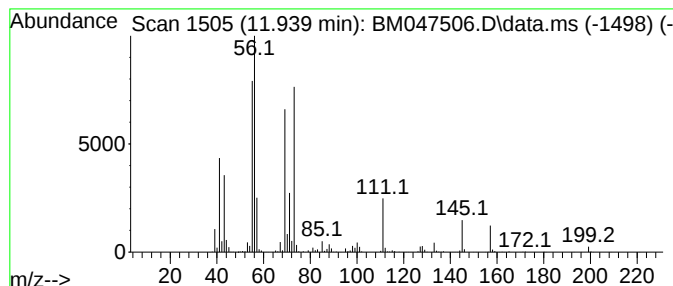
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 unknown-02 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.939	5.07 ng/ul	3318290	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Neopentyl glycol	104	C5H12O2	000126-30-7	33
2		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	32
3		1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	32
4		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	32
5		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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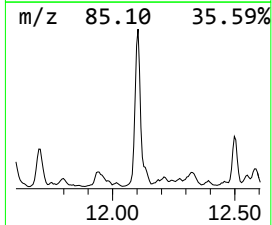
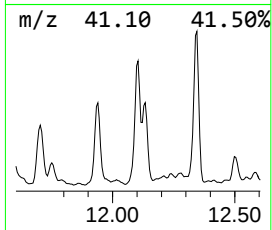
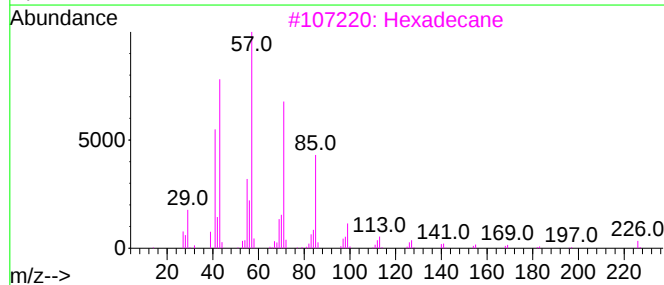
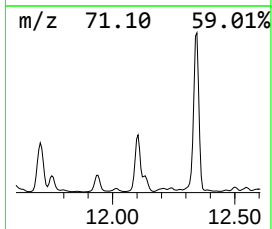
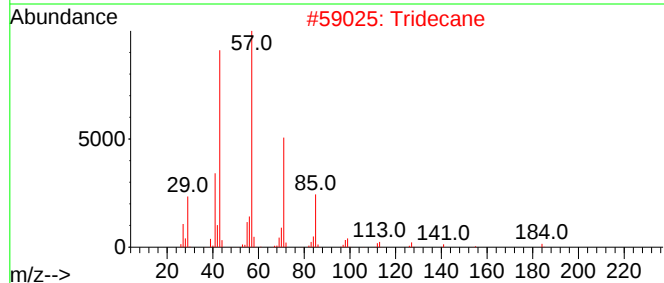
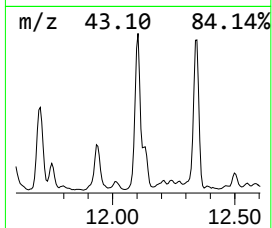
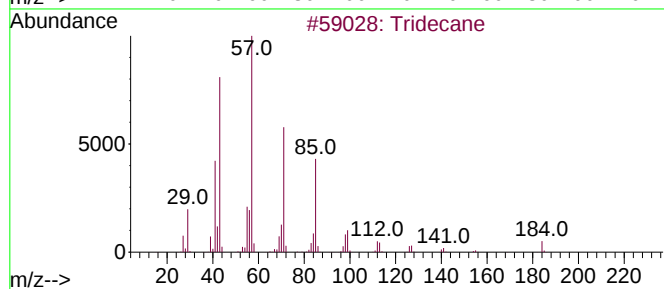
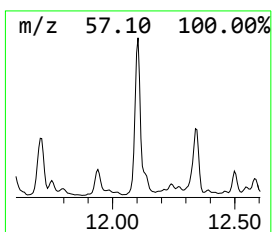
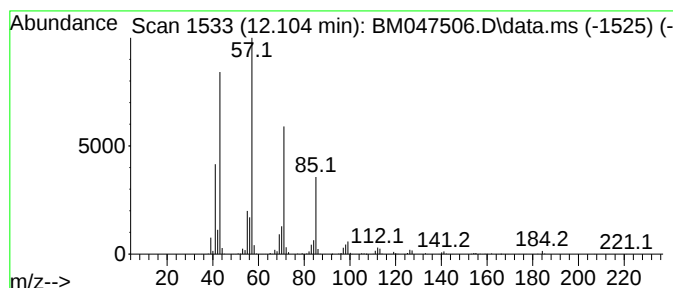
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	5.93 ng/ul	3884420	Naphthalene-d8	10.663

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	93
2		Tridecane	184	C13H28	000629-50-5	93
3		Hexadecane	226	C16H34	000544-76-3	91
4		Hexadecane	226	C16H34	000544-76-3	90
5		Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

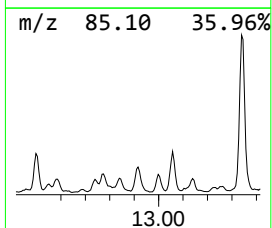
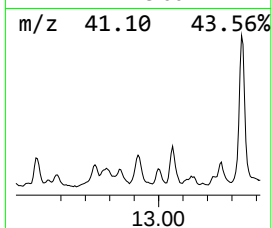
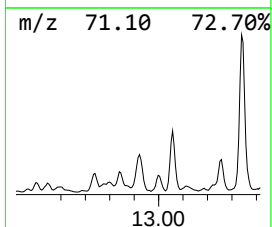
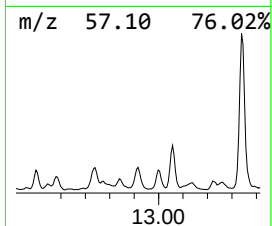
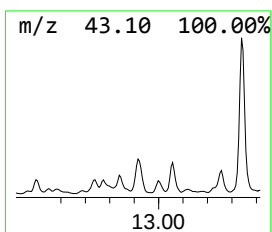
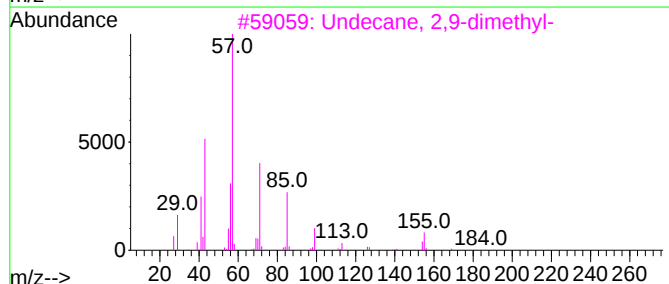
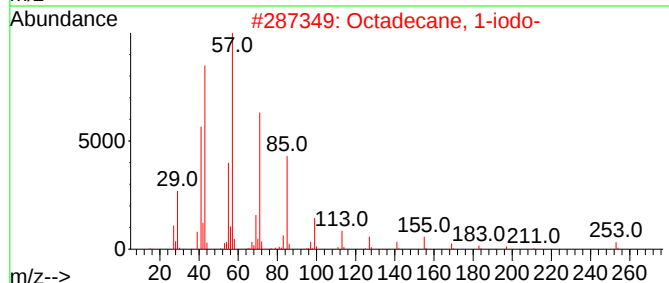
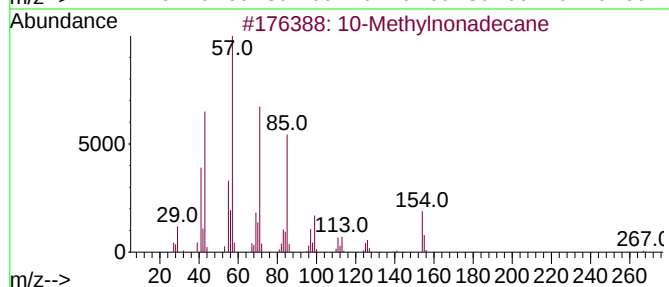
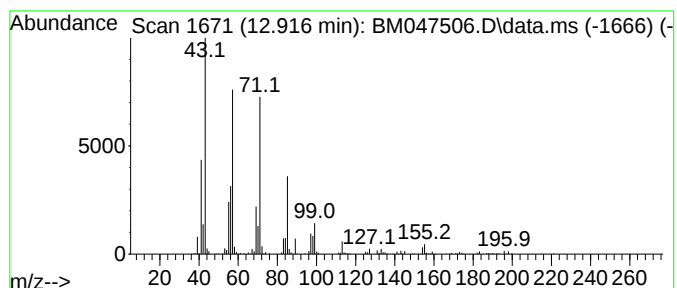
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.916	3.80 ng/ul	1246750	Acenaphthene-d10	14.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane	282	C20H42	056862-62-5	72
2	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64
3	Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	59
4	Dodecane	170	C12H26	000112-40-3	59
5	Sulfurous acid, pentyl undecyl e...	306	C16H34O3S	1000309-14-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

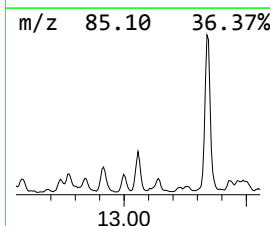
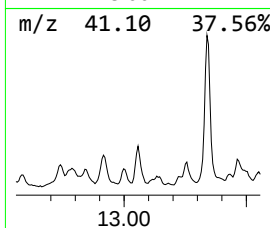
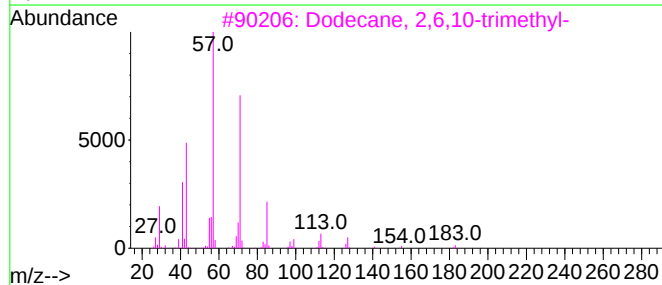
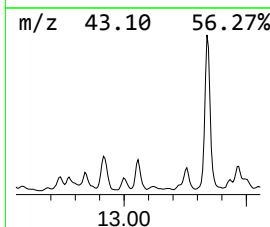
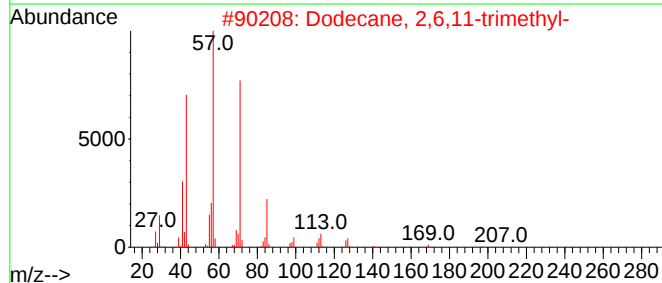
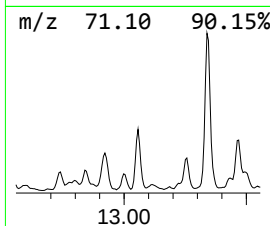
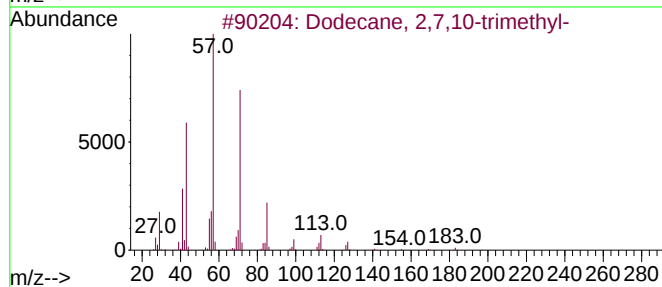
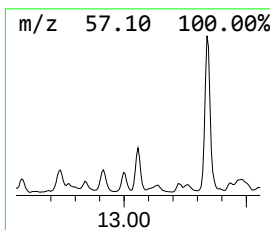
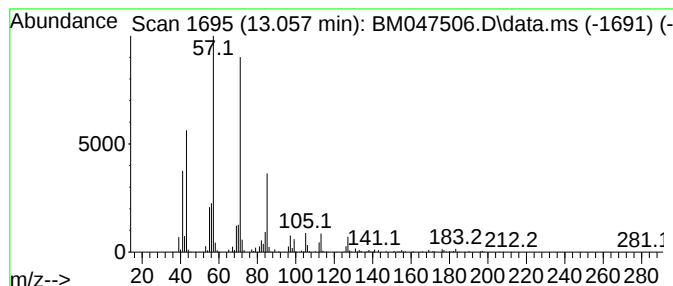
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.057	3.93 ng/ul	1290770	Acenaphthene-d10			14.492
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	86
2	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	78
3	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	72
4	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	60
5	Dodecane, 4,6-dimethyl-		198	C14H30	061141-72-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
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Sample : P3878-08DL2 20X
Misc :
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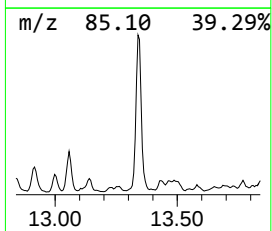
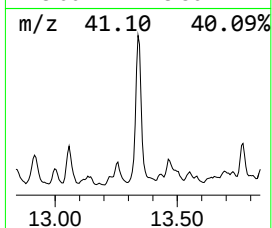
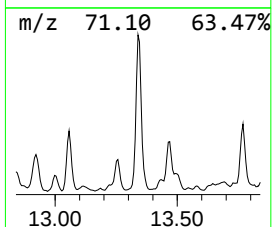
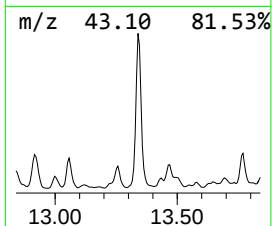
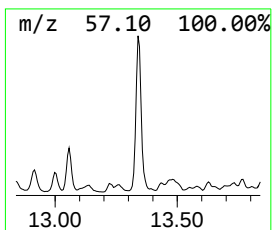
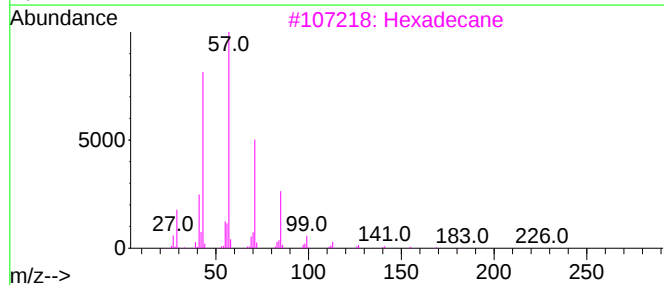
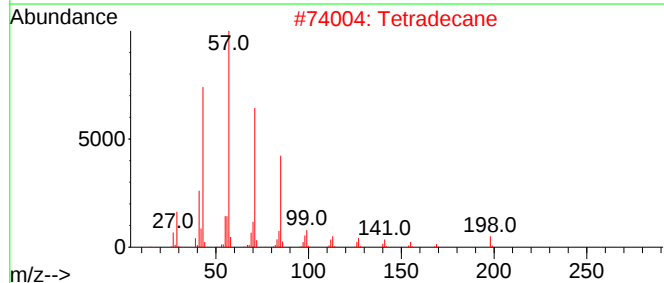
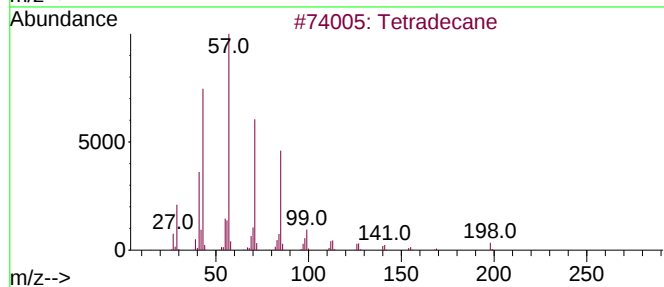
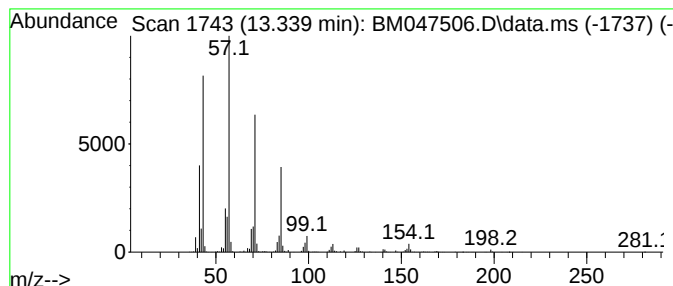
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.339	14.10 ng/ul	4628380	Acenaphthene-d10	14.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	95
2	Tetradecane	198	C14H30	000629-59-4	93
3	Hexadecane	226	C16H34	000544-76-3	91
4	Hexadecane	226	C16H34	000544-76-3	91
5	Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

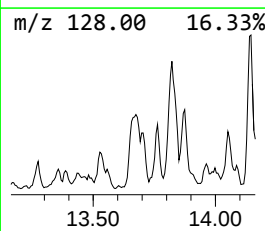
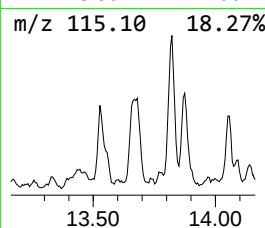
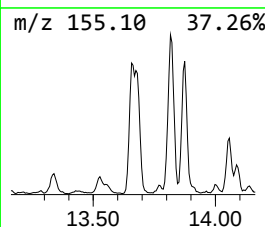
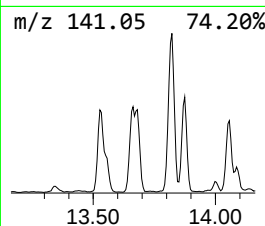
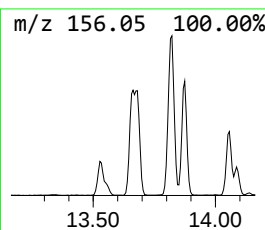
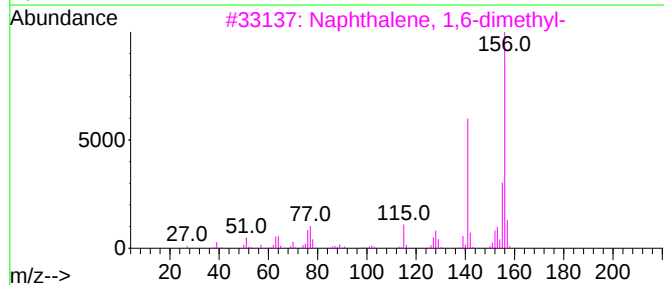
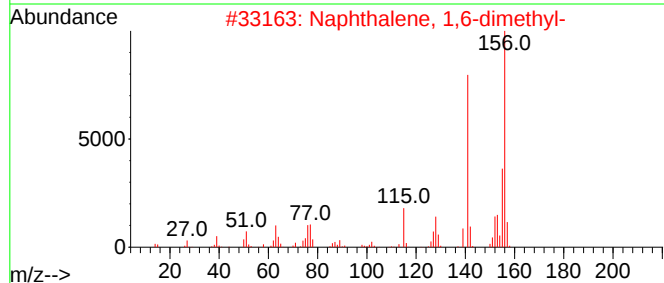
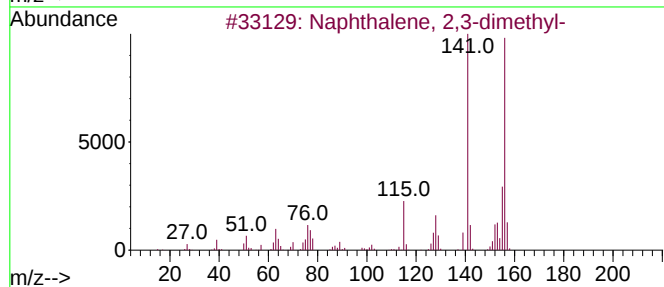
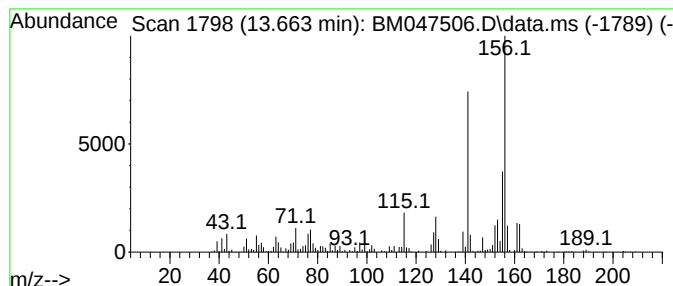
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 Naphthalene, 2,3-dimethyl- Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	4.02 ng/ul	1320590	Acenaphthene-d10	14.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

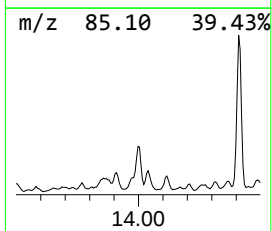
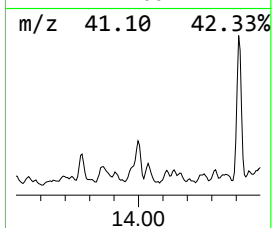
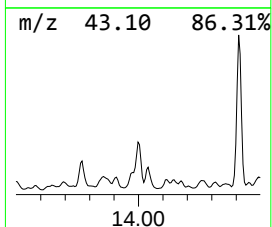
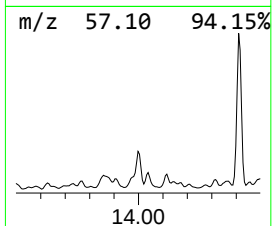
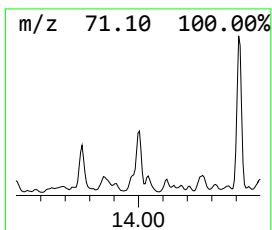
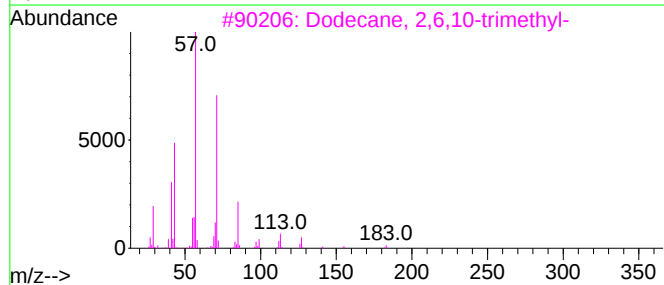
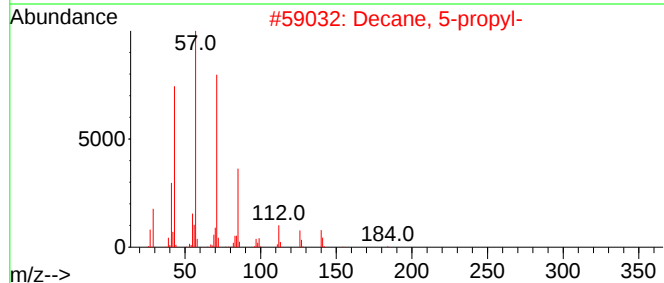
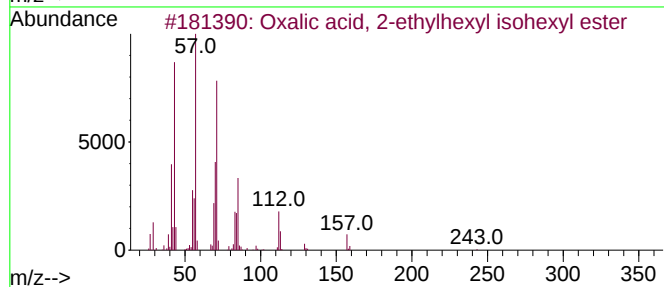
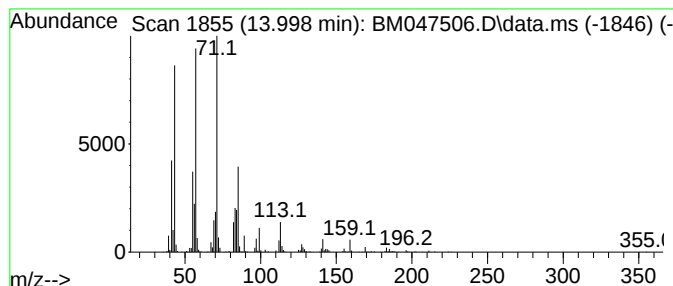
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 Oxalic acid, 2-ethylhexyl i... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.998	7.28 ng/ul	2389860	Acenaphthene-d10	14.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl isohex...	286	C16H30O4	1000309-38-8	86
2			Decane, 5-propyl-	184	C13H28	017312-62-8	62
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	58
4			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	53
5			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

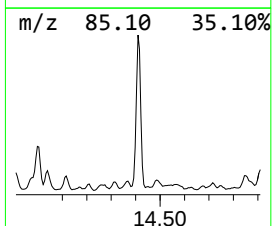
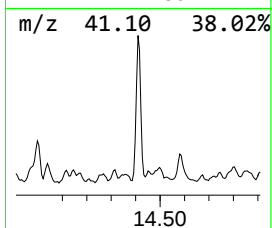
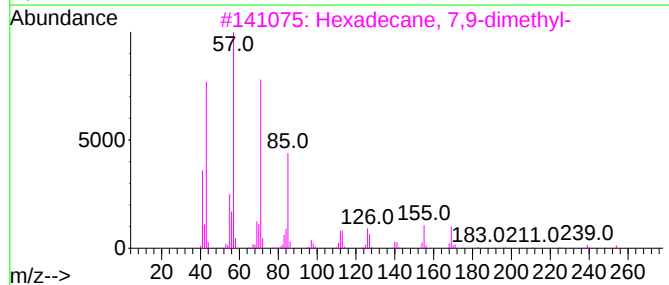
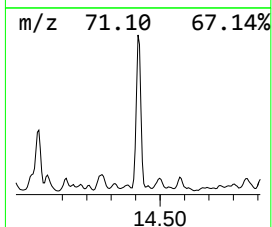
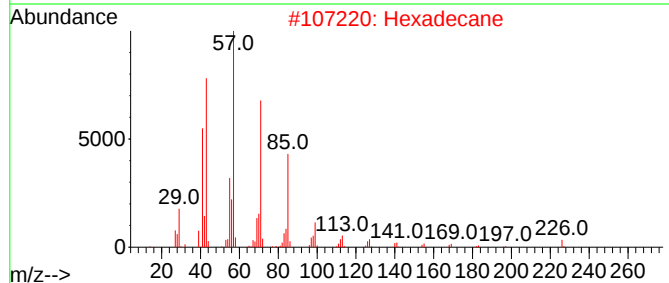
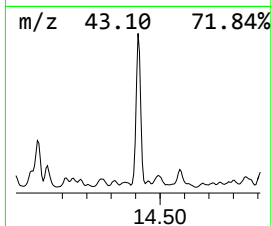
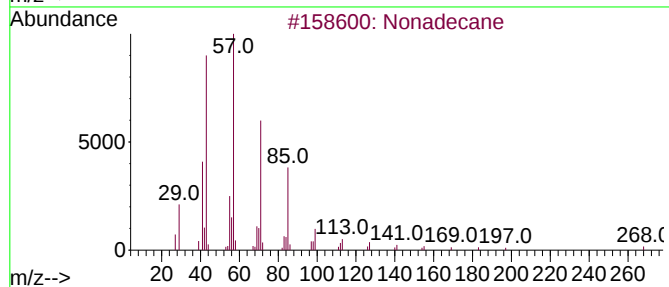
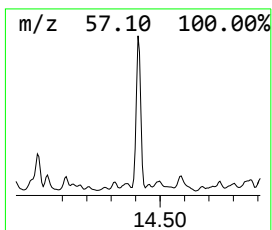
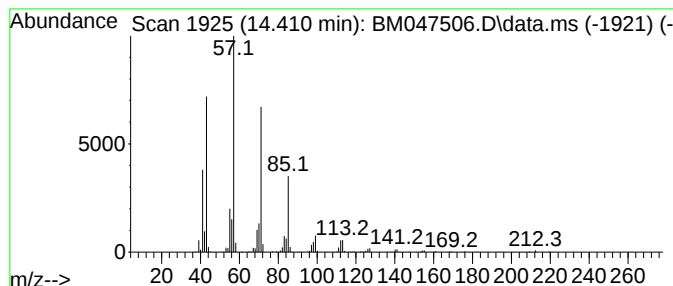
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.410	14.48 ng/ul	4752840	Acenaphthene-d10	14.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	91
2	Hexadecane	226	C16H34	000544-76-3	86
3	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	86
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86
5	Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

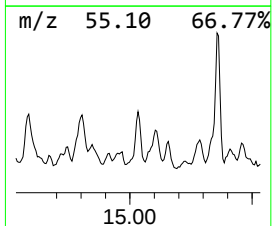
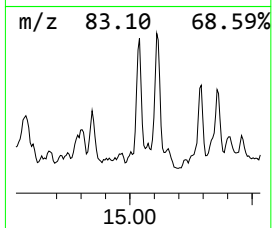
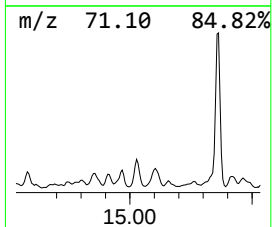
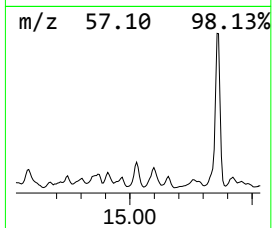
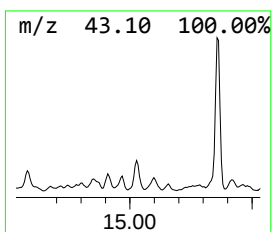
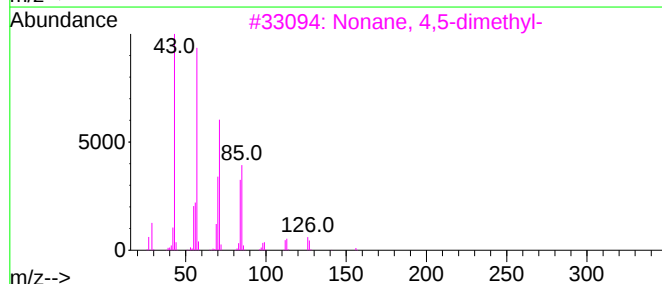
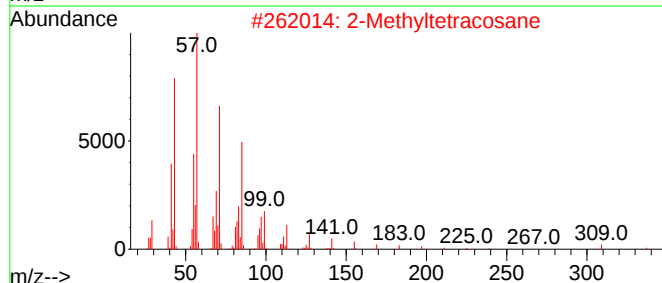
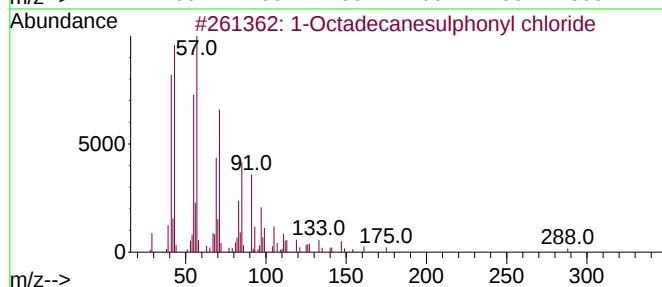
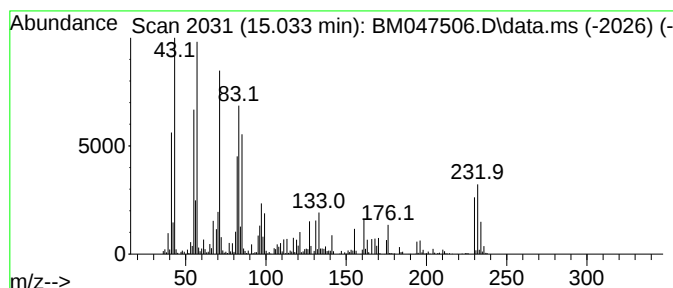
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 unknown-03 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.033	4.66 ng/ul	1529330	Acenaphthene-d10	14.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	46
2	2-Methyltetracosane	352	C25H52	001560-78-7	46
3	Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	42
4	10-Methylnonadecane	282	C20H42	056862-62-5	38
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

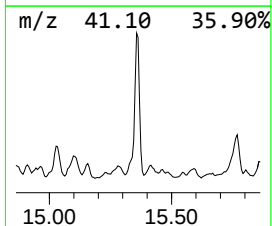
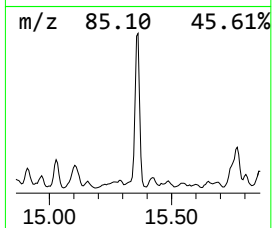
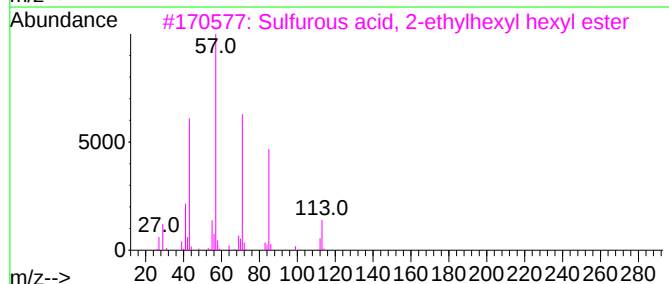
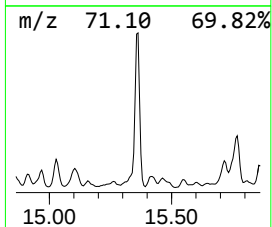
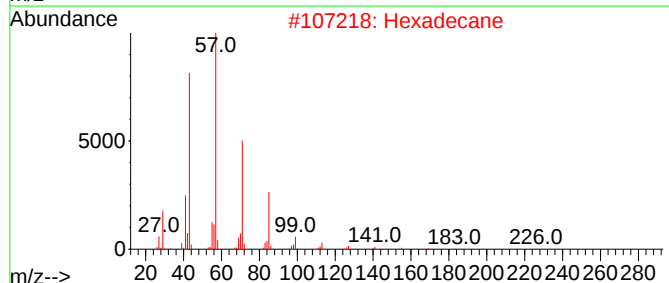
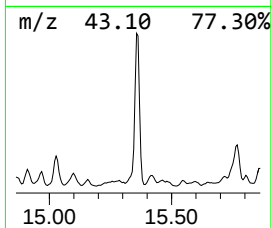
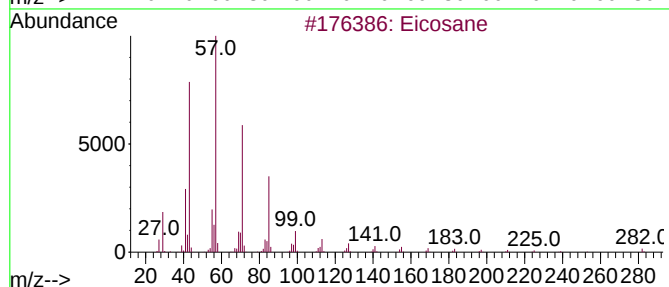
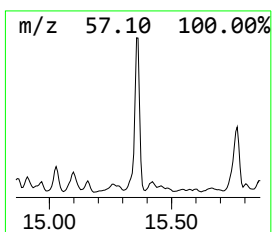
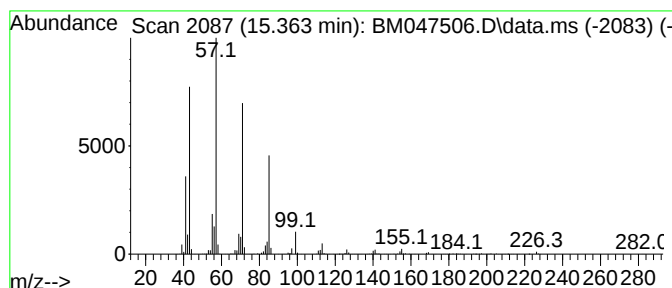
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.363	12.29 ng/ul	4033400	Acenaphthene-d10		14.492	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Hexadecane		226	C16H34	000544-76-3	80
3	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	78
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	72
5	Decane, 3-bromo-		220	C10H21Br	030571-71-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

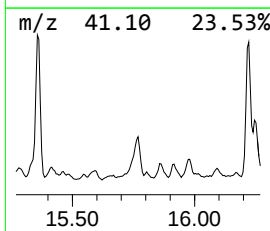
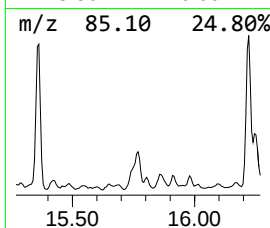
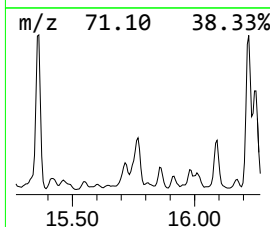
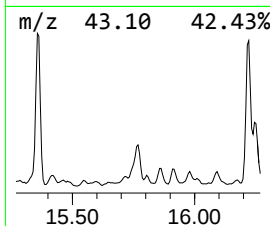
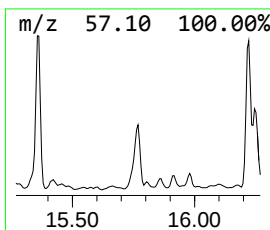
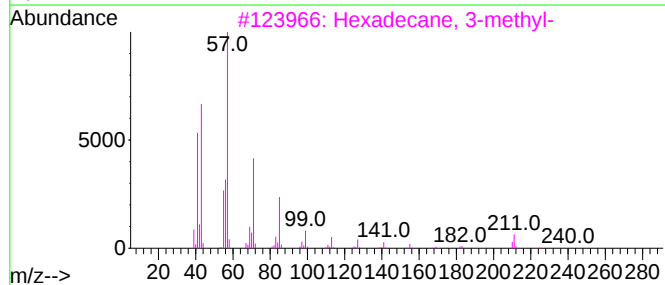
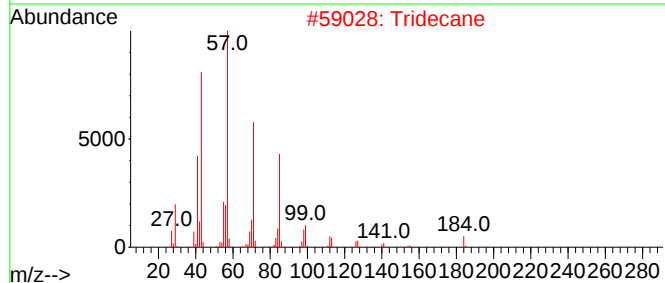
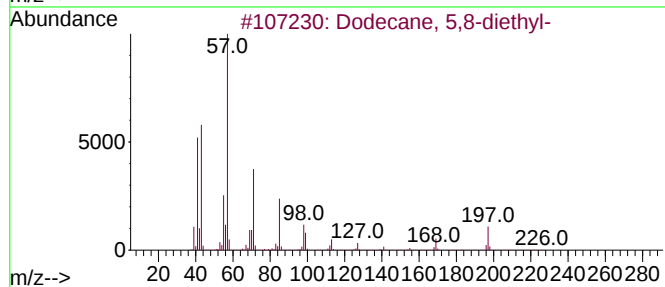
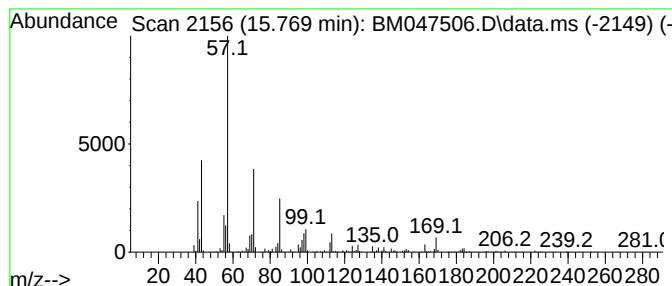
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.769	6.41 ng/ul	2102550	Acenaphthene-d10		14.492	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 5,8-diethyl-		226	C16H34	024251-86-3	87
2	Tridecane		184	C13H28	000629-50-5	81
3	Hexadecane, 3-methyl-		240	C17H36	006418-43-5	80
4	3,5-Dimethyldodecane		198	C14H30	107770-99-0	80
5	Decane, 2,6,8-trimethyl-		184	C13H28	062108-26-3	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

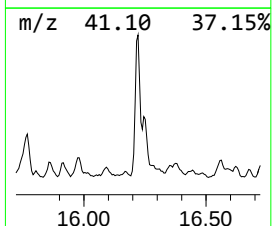
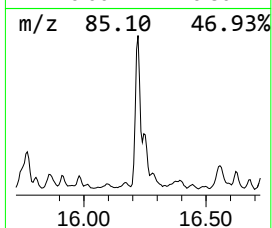
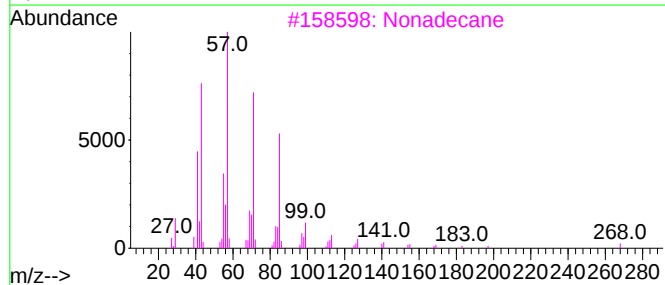
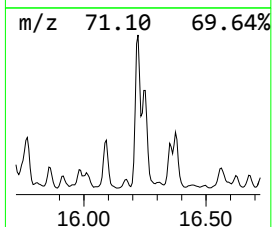
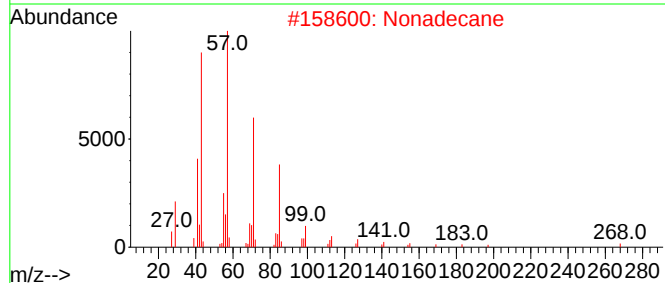
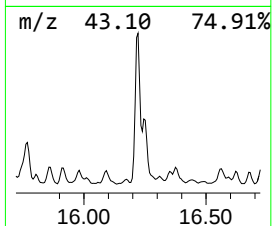
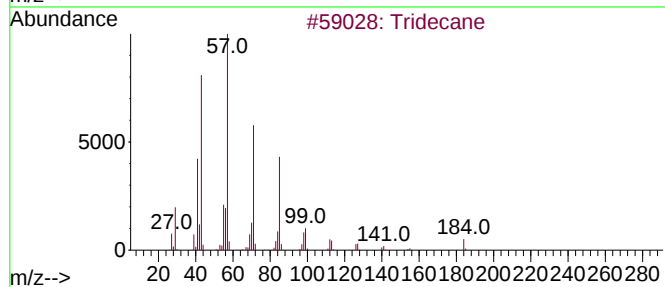
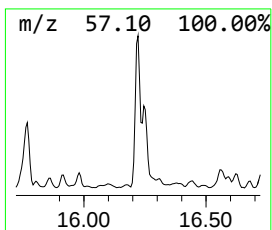
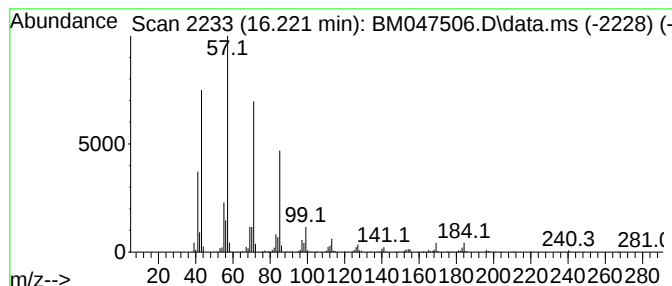
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.221	16.03 ng/ul	4944540	Phenanthrene-d10	17.233	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Nonadecane	268	C19H40	000629-92-5	91
3	Nonadecane	268	C19H40	000629-92-5	90
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5	Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

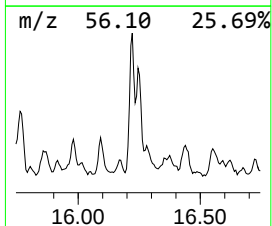
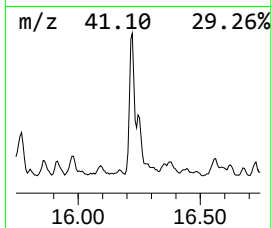
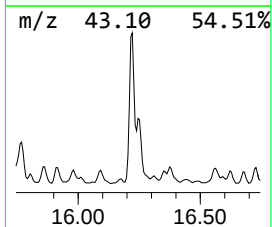
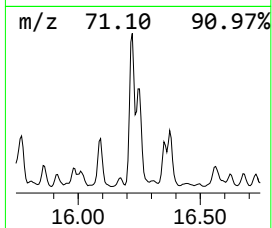
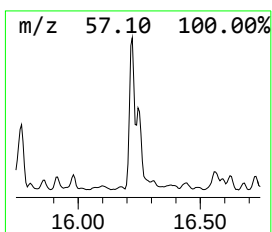
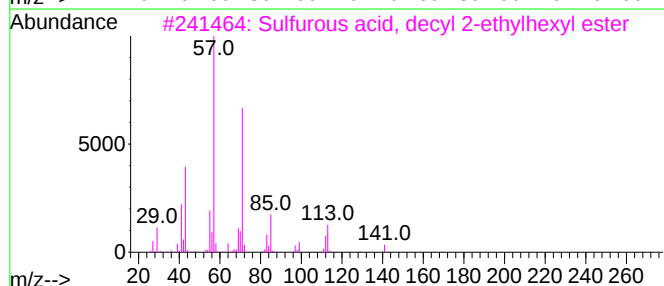
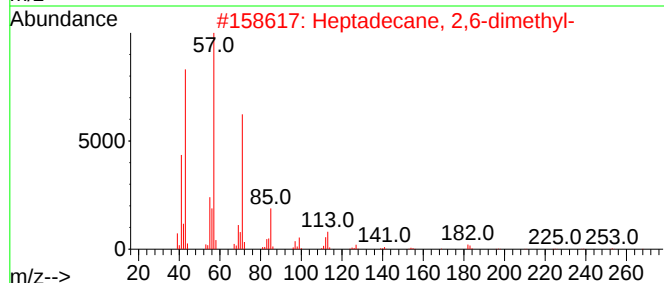
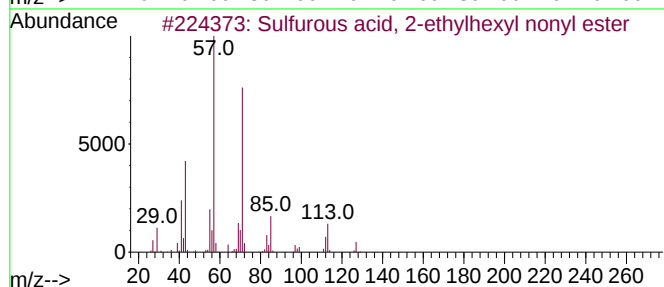
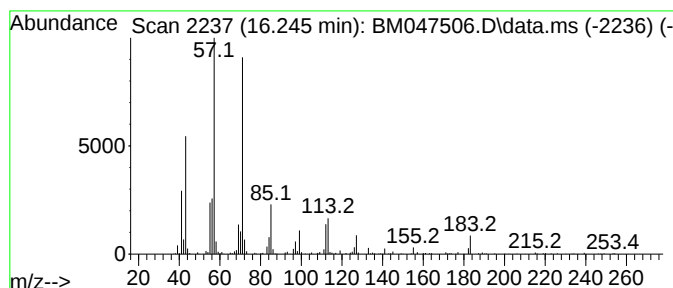
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 Sulfurous acid, 2-ethylhexy... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.245	5.10 ng/ul	1572590	Phenanthrene-d10	17.233	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	83
2	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	83
3	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	80
4	Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	72
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

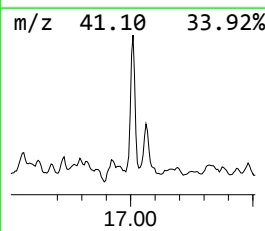
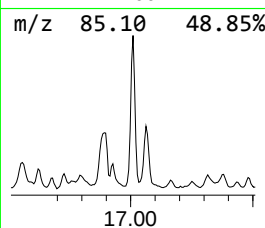
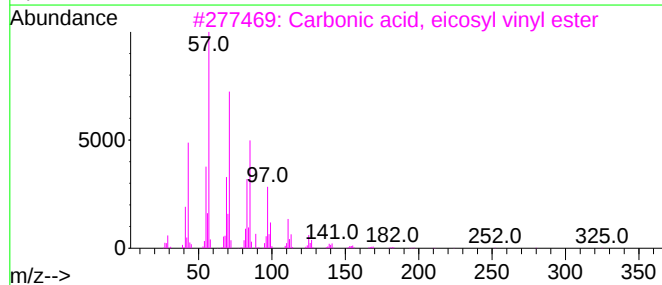
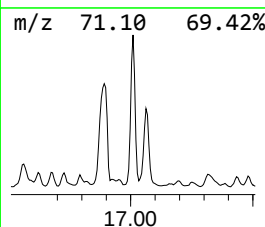
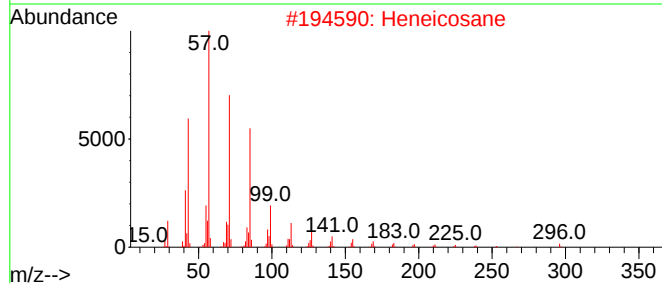
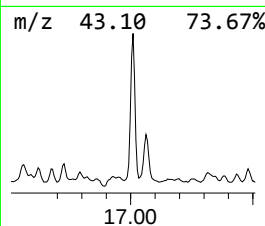
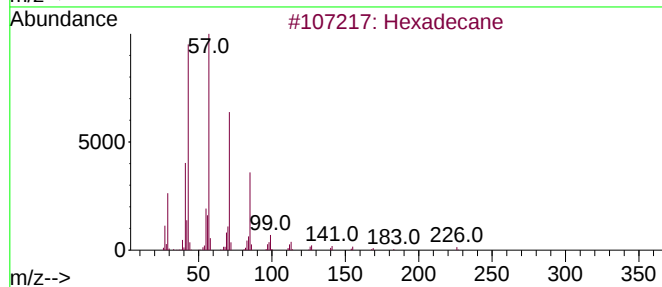
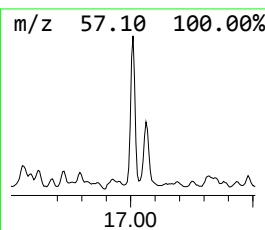
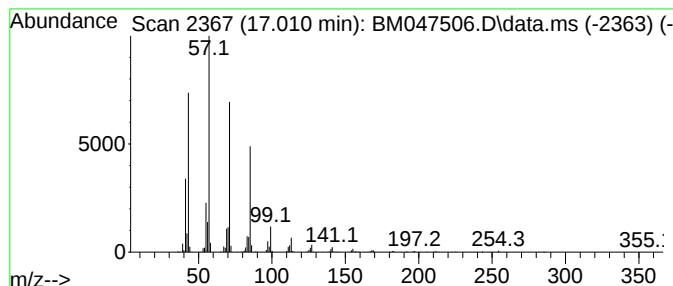
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.010	10.10 ng/ul	3114290	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	90
2			Heneicosane	296	C21H44	000629-94-7	90
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	87
4			Heptadecane	240	C17H36	000629-78-7	87
5			Tridecane, 6-propyl-	226	C16H34	055045-10-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

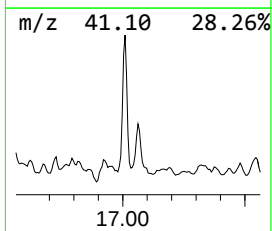
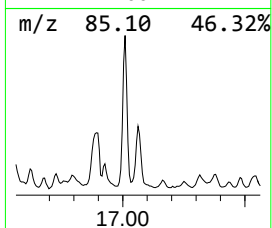
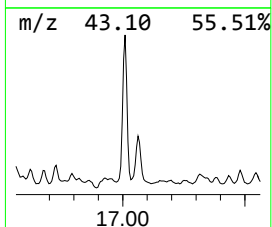
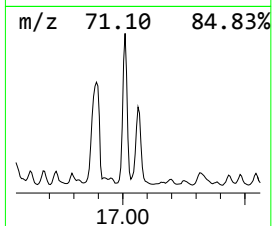
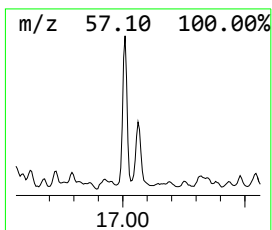
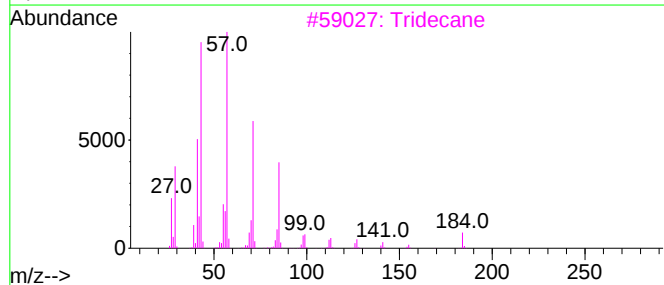
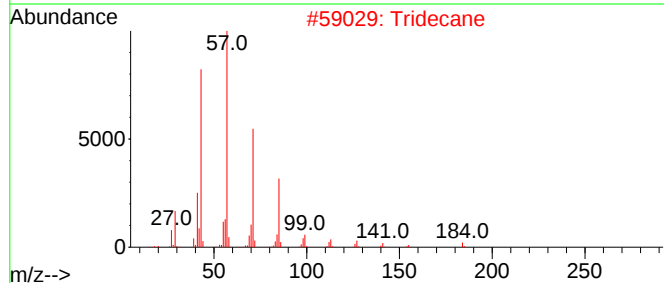
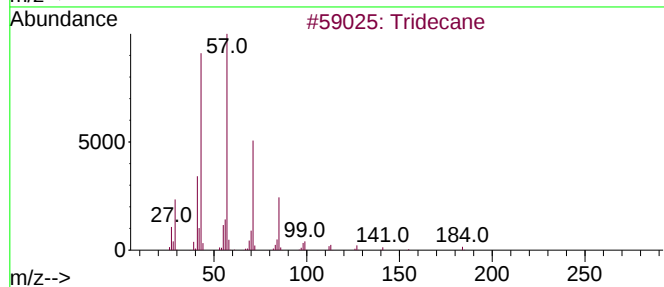
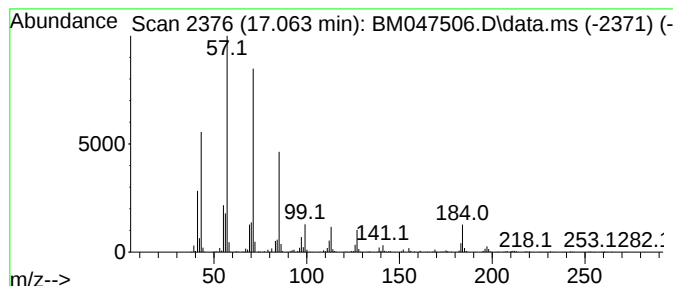
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.063	7.48 ng/ul	2308650	Phenanthrene-d10	17.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	91
2			Tridecane	184	C13H28	000629-50-5	91
3			Tridecane	184	C13H28	000629-50-5	87
4			Tetracosane	338	C24H50	000646-31-1	86
5			Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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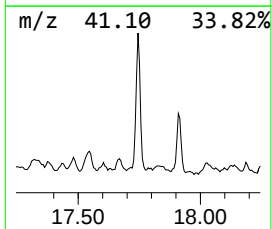
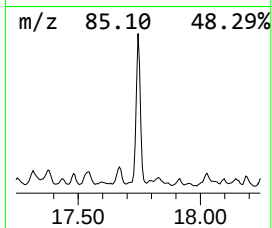
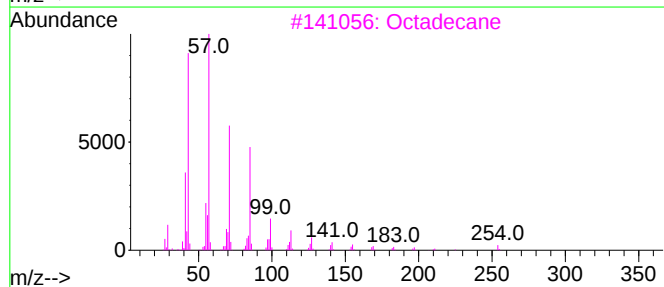
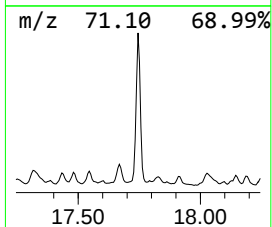
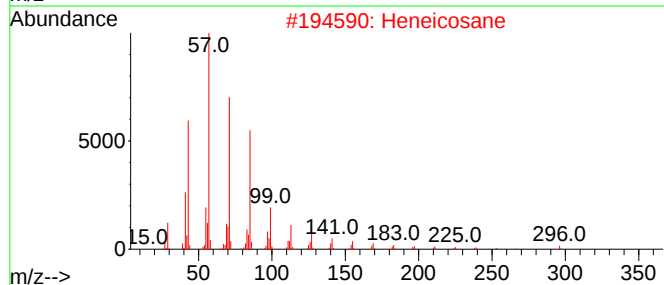
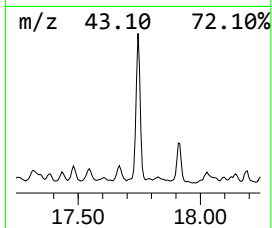
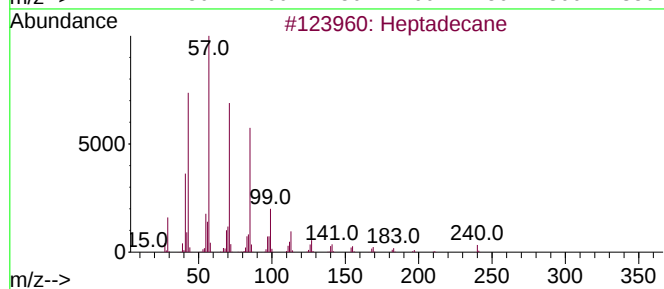
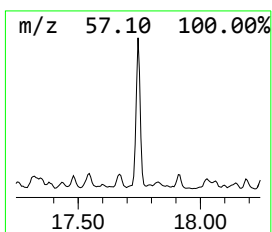
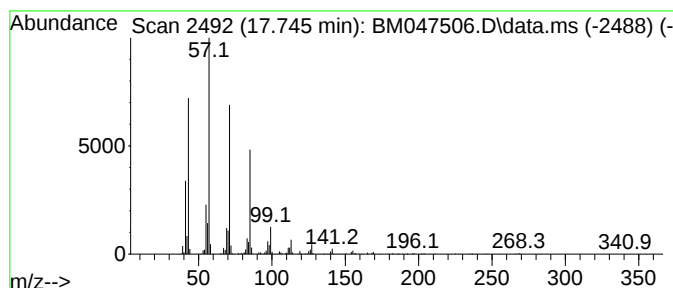
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.745	9.55 ng/ul	2946920	Phenanthrene-d10	17.233

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octadecane	254	C18H38	000593-45-3	90
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

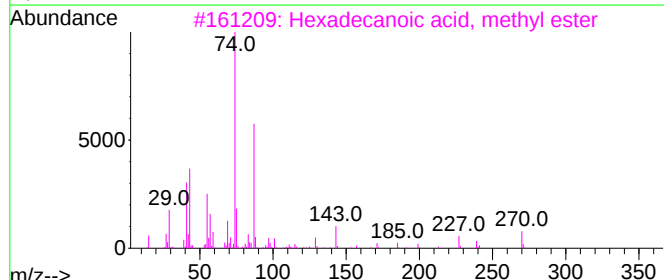
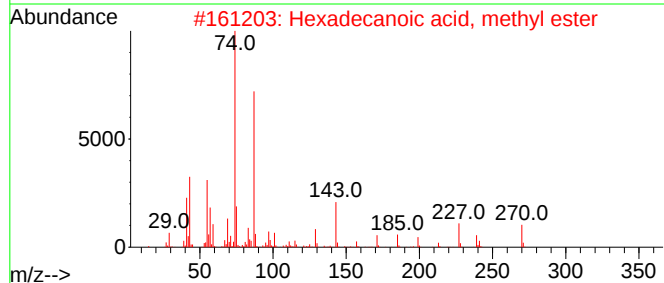
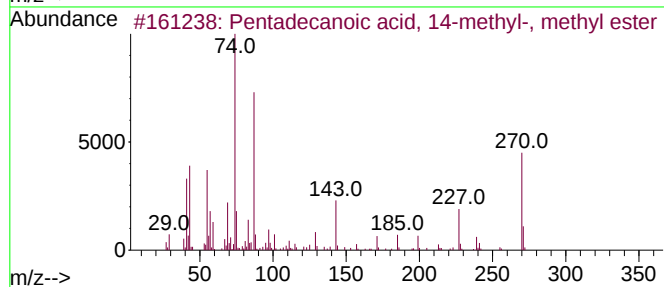
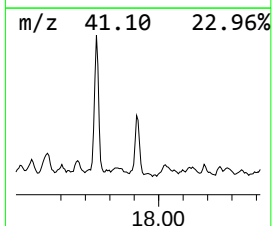
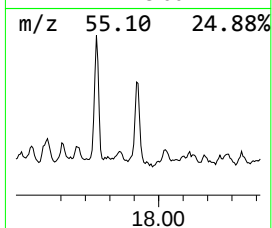
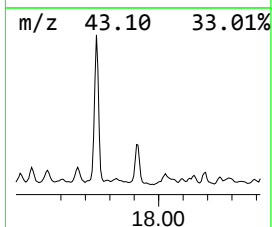
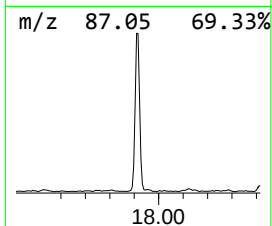
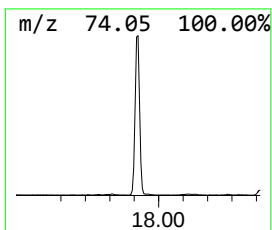
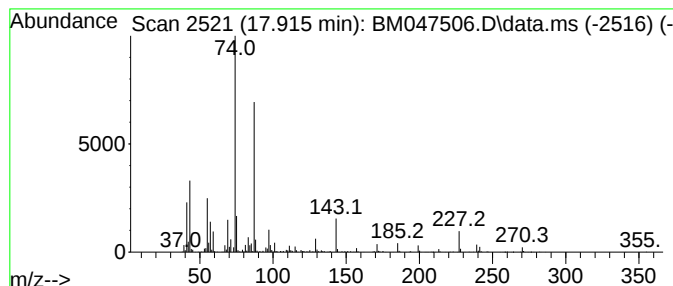
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 69 Pentadecanoic acid, 14-meth... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.915	5.30 ng/ul	1634470	Phenanthrene-d10	17.233	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	92
3	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	87
5	Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

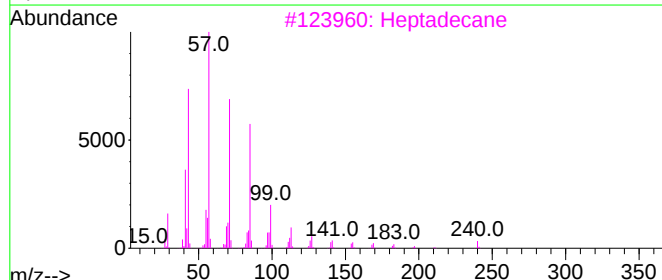
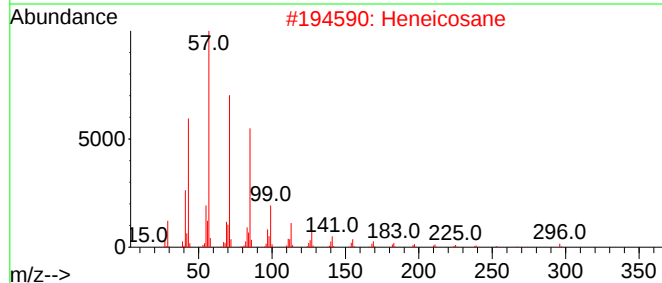
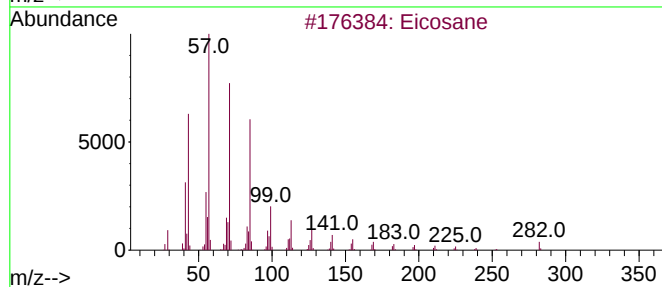
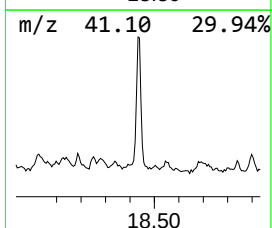
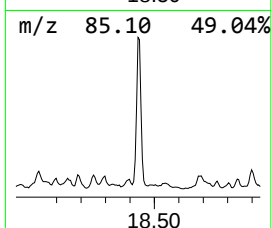
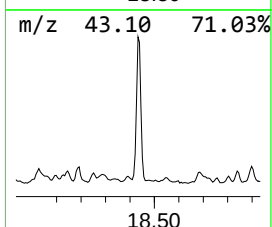
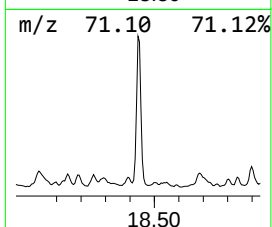
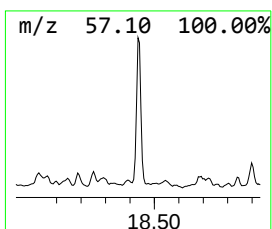
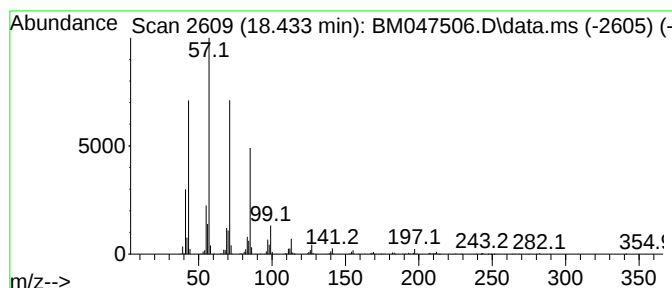
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.433	7.40 ng/ul	2282450	Phenanthrene-d10	17.233

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	97
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane	240	C17H36	000629-78-7	91
4		Nonadecane	268	C19H40	000629-92-5	90
5		Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

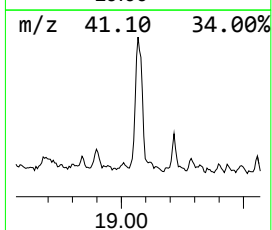
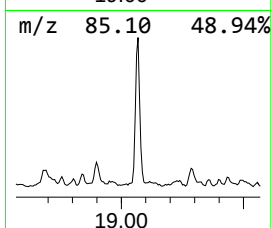
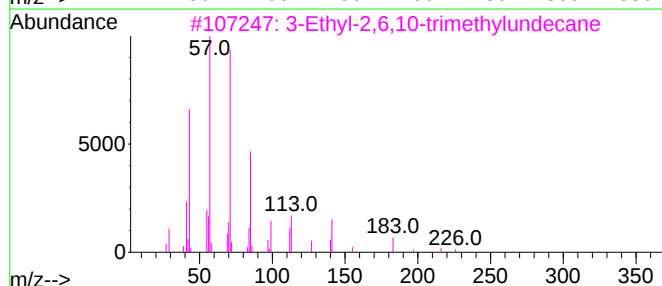
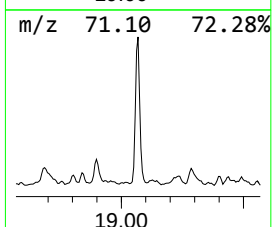
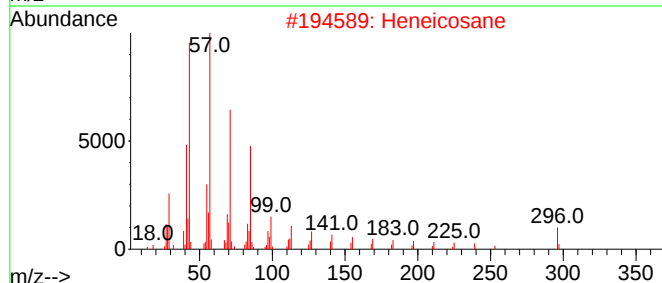
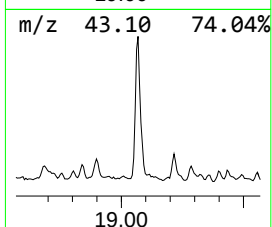
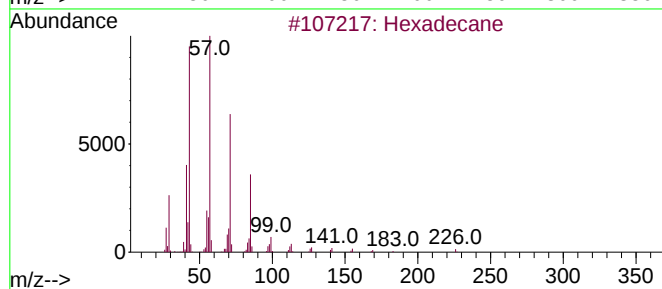
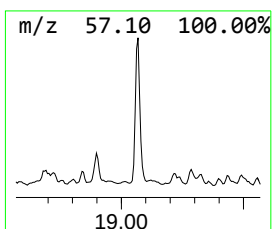
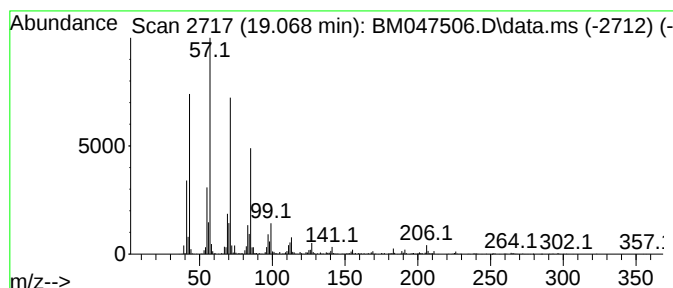
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.068	10.70 ng/ul	3300240	Phenanthrene-d10		17.233	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	95
2	Heneicosane		296	C21H44	000629-94-7	94
3	3-Ethyl-2,6,10-trimethylundecane		226	C16H34	1000432-25-9	93
4	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	93
5	Hexacosane		366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

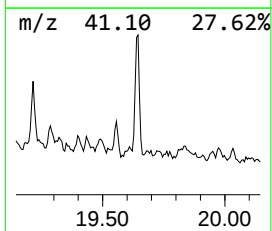
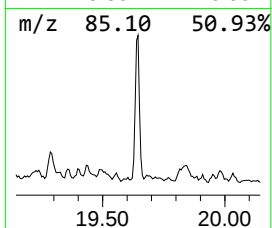
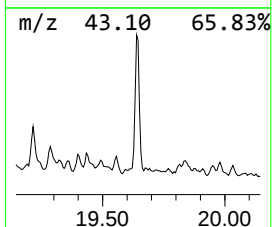
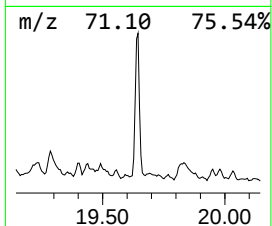
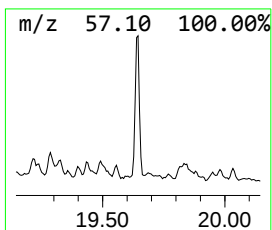
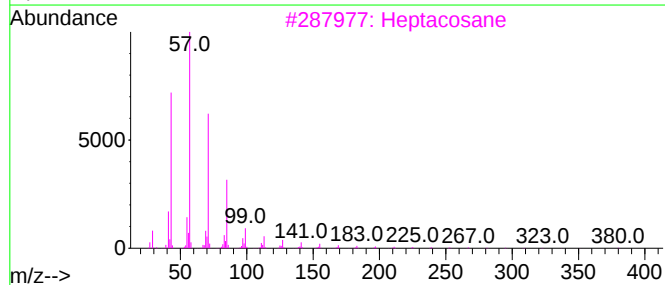
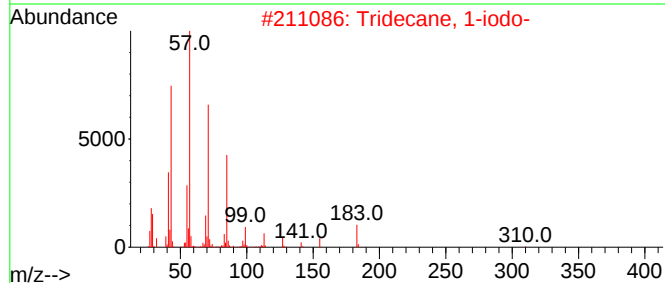
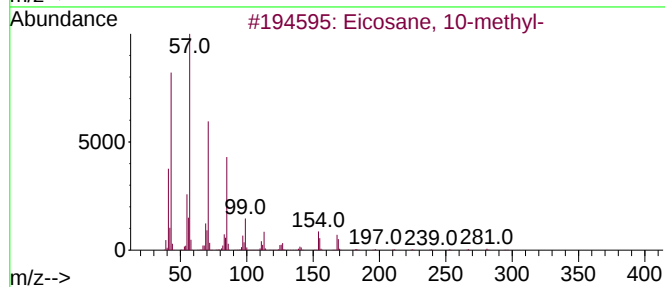
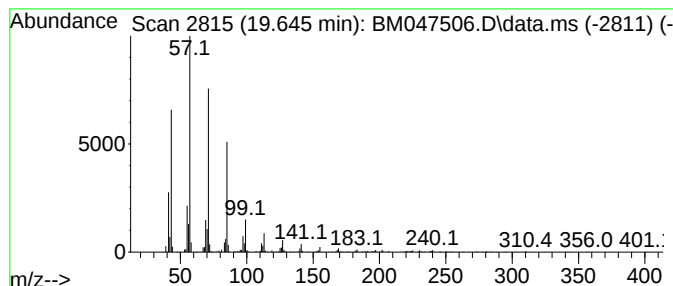
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 73 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.645	4.07 ng/ul	1264680	Chrysene-d12	21.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	94
3	Heptacosane	380	C27H56	000593-49-7	91
4	Hexadecane	226	C16H34	000544-76-3	91
5	Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

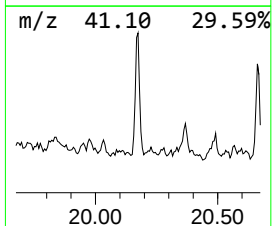
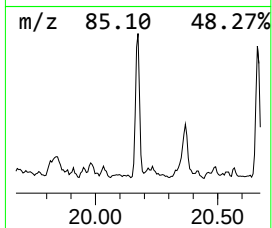
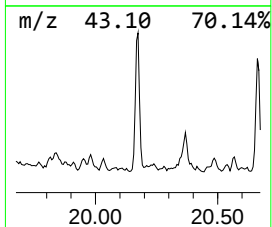
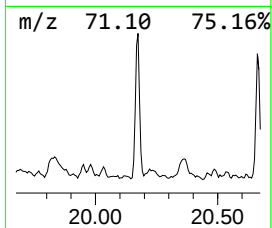
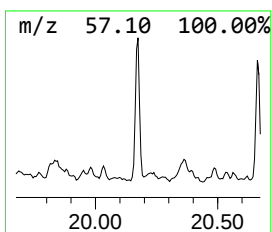
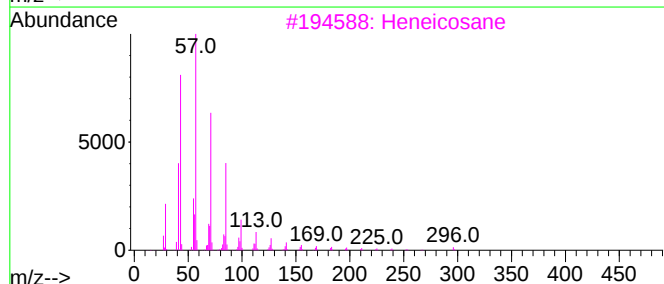
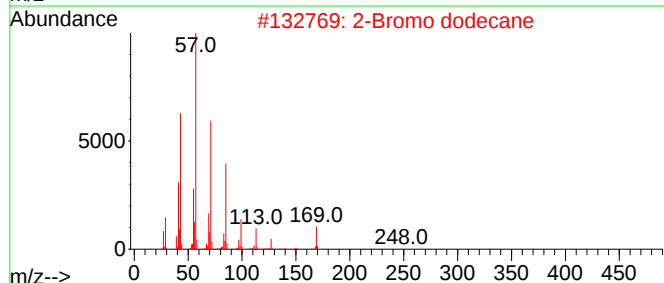
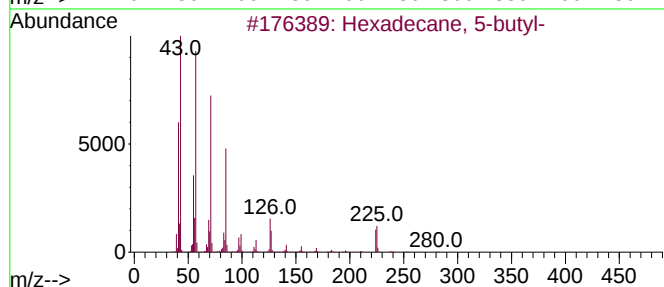
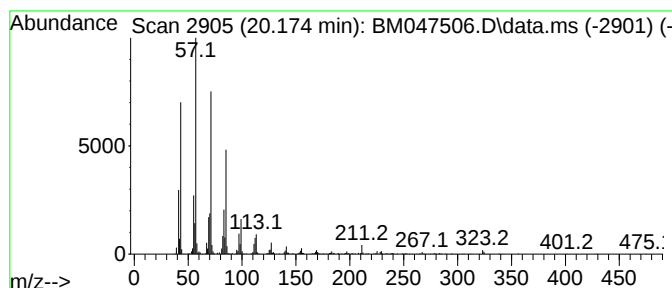
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.174	3.55 ng/ul	1104980	Chrysene-d12	21.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	91
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
3	Heneicosane	296	C21H44	000629-94-7	87
4	Hexadecane	226	C16H34	000544-76-3	80
5	Tridecane, 7-propyl-	226	C16H34	055045-09-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

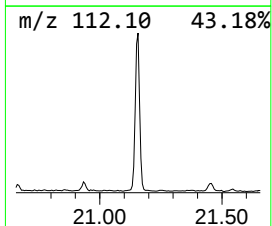
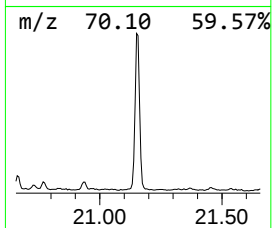
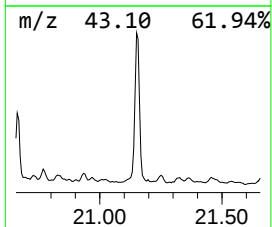
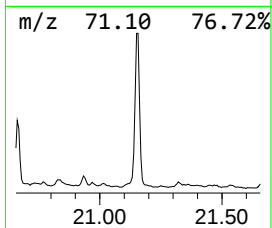
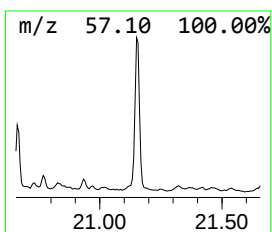
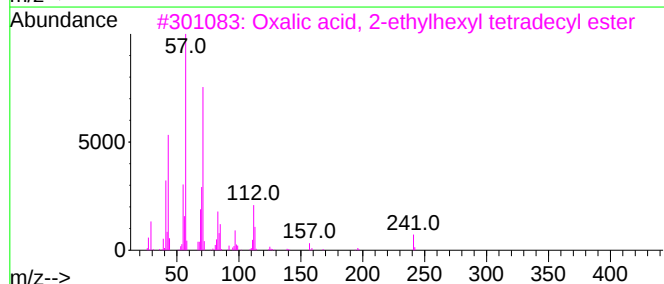
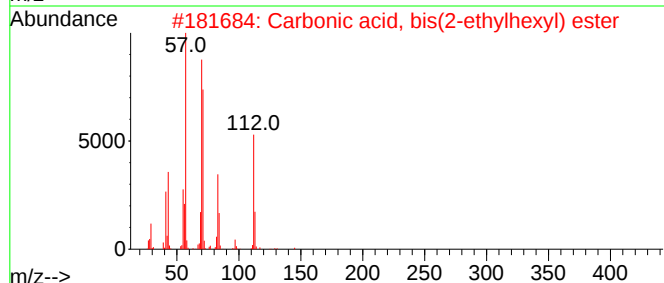
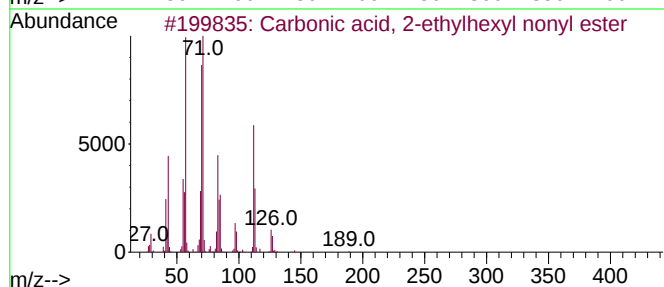
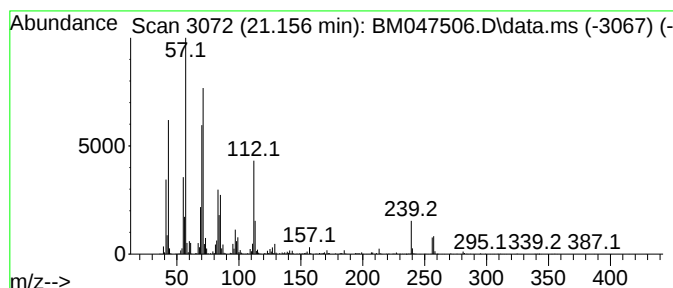
Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 Carbonic acid, 2-ethylhexyl... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.156	9.91 ng/ul	3082240	Chrysene-d12	21.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, 2-ethylhexyl nonyl...	300	C18H36O3	1000383-13-6	80
2	Carbonic acid, bis(2-ethylhexyl)...	286	C17H34O3	014858-73-2	58
3	Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	58
4	3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	49
5	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

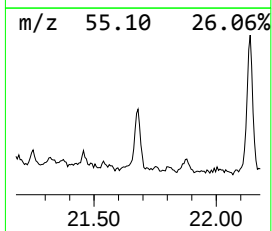
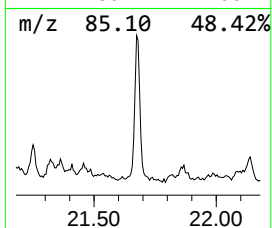
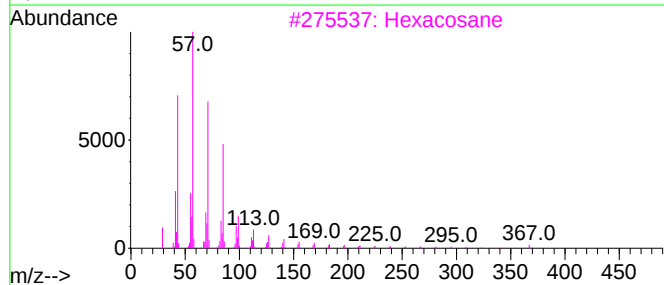
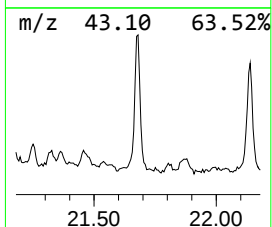
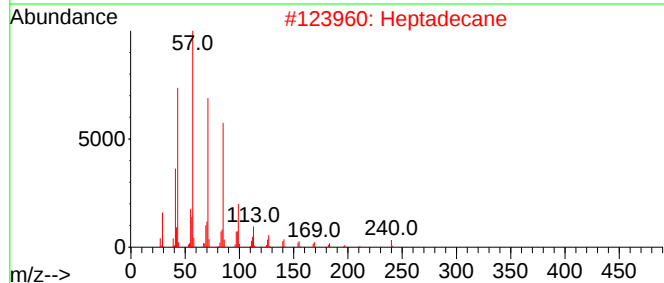
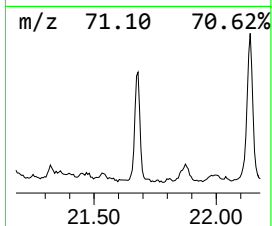
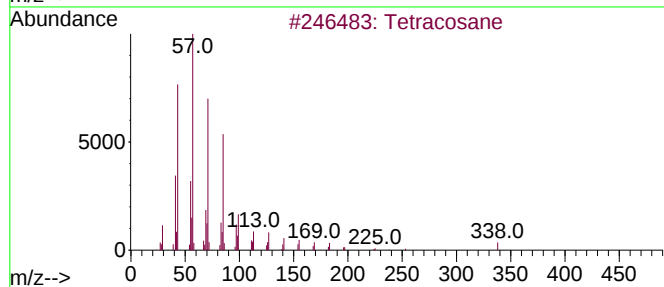
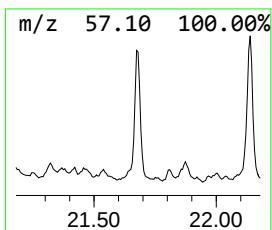
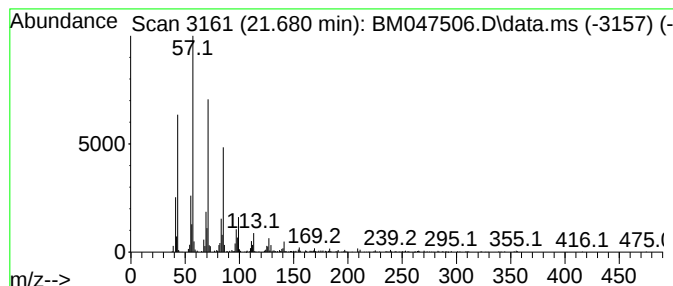
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 72

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.680	2.83 ng/ul	879826	Chrysene-d12		21.456	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	91
2	Heptadecane		240	C17H36	000629-78-7	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	91
5	Heptacosane		380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

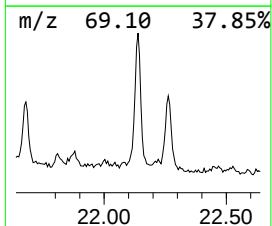
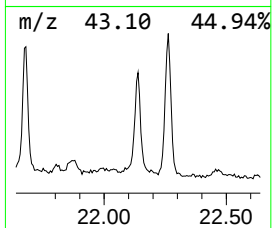
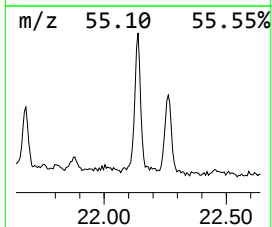
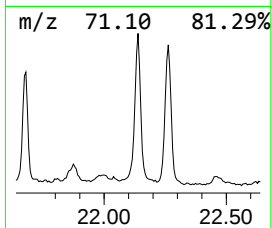
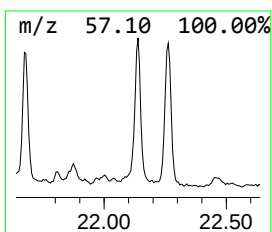
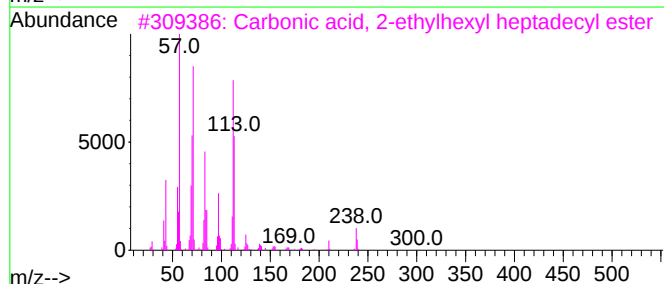
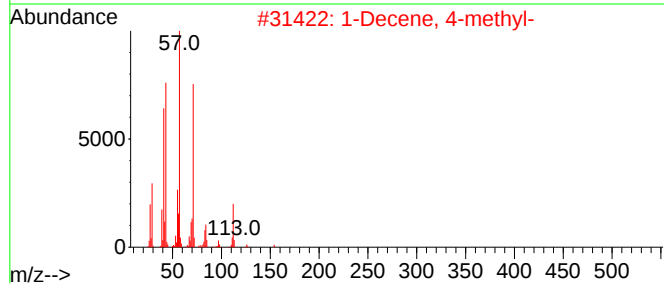
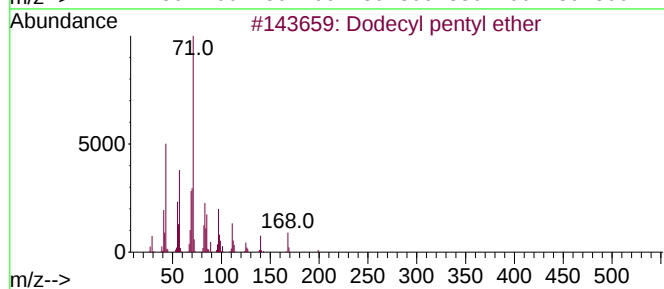
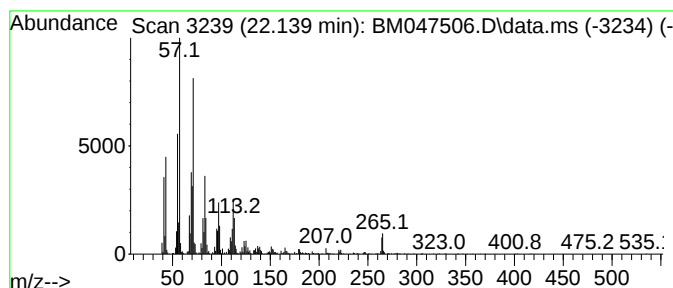
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 77 Dodecyl pentyl ether Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.139	5.06 ng/ul	1575440	Chrysene-d12	21.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecyl pentyl ether	256	C17H36O	1010406-41-6	53
2	1-Decene, 4-methyl-	154	C11H22	013151-29-6	53
3	Carbonic acid, 2-ethylhexyl hept...	412	C26H52O3	1000383-14-4	53
4	Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	49
5	Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
Data File : BM047506.D
Acq On : 12 Sep 2024 16:49
Operator : RC/JU
Sample : P3878-08DL2 20X
Misc :
ALS Vial : 11 Sample Multiplier: 1

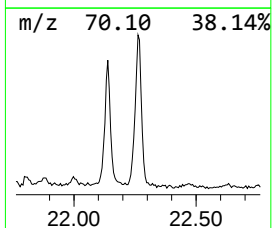
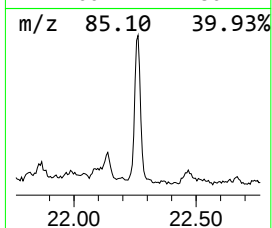
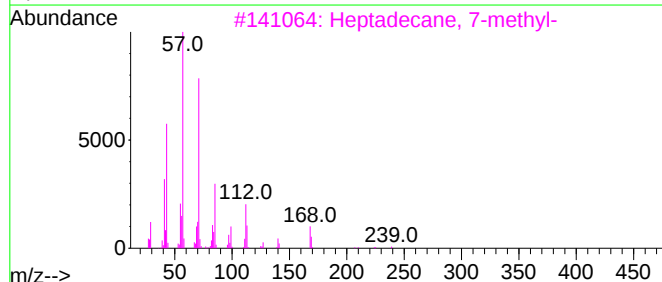
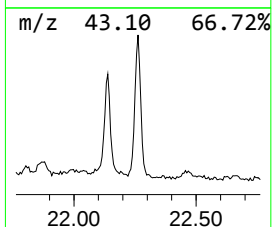
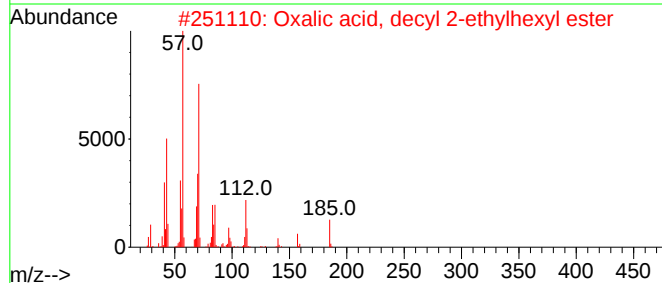
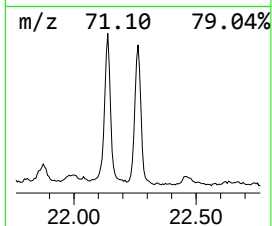
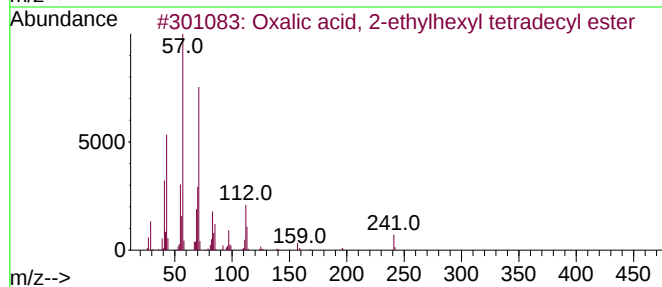
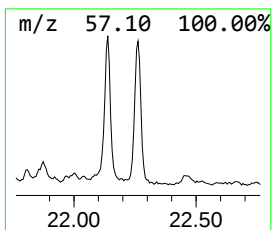
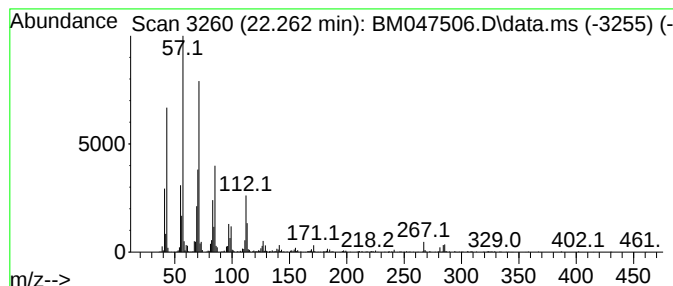
Instrument :
BNA_M
ClientSampleId :
DCVY6DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 78 Oxalic acid, 2-ethylhexyl t... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.262	4.17 ng/ul	1296330	Chrysene-d12	21.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	83
2	Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	80
3	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	72
4	Tetratetracontane	619	C44H90	007098-22-8	64
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
 Data File : BM047506.D
 Acq On : 12 Sep 2024 16:49
 Operator : RC/JU
 Sample : P3878-08DL2 20X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVY6DL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butanoic acid	3.999	9.7	ng/u1	2485660	1	7.869	5101570	20.0
(DEL) Alkane: S...	5.899	14.5	ng/u1	3700160	1	7.869	5101570	20.0
Hexylene glycol	6.175	11.3	ng/u1	2875480	1	7.869	5101570	20.0
unknown-01	6.369	6.1	ng/u1	1550690	1	7.869	5101570	20.0
(DEL) Alkane: S...	6.440	4.8	ng/u1	1233490	1	7.869	5101570	20.0
Cyclohexanone, ...	6.516	5.3	ng/u1	1342740	1	7.869	5101570	20.0
(DEL) Alkane: S...	6.551	3.4	ng/u1	878358	1	7.869	5101570	20.0
(DEL) Alkane: S...	6.781	4.0	ng/u1	1024620	1	7.869	5101570	20.0
Benzene, propyl-	6.846	3.9	ng/u1	990243	1	7.869	5101570	20.0
(DEL) Alkane: S...	6.875	7.4	ng/u1	1891570	1	7.869	5101570	20.0
(DEL) Alkane: S...	6.934	9.5	ng/u1	2433930	1	7.869	5101570	20.0
Benzene, 1-ethy...	6.963	6.1	ng/u1	1559760	1	7.869	5101570	20.0
(DEL) Alkane: S...	7.040	15.1	ng/u1	3838700	1	7.869	5101570	20.0
Mesitylene	7.098	4.7	ng/u1	1188420	1	7.869	5101570	20.0
Benzene, 1-ethy...	7.263	5.9	ng/u1	1508730	1	7.869	5101570	20.0
Carbonic acid, ...	7.298	5.7	ng/u1	1460870	1	7.869	5101570	20.0
Butanoic acid, ...	7.393	7.9	ng/u1	2003730	1	7.869	5101570	20.0
(DEL) Alkane: S...	7.510	59.7	ng/u1	15238500	1	7.869	5101570	20.0
(DEL) Alkane: S...	7.716	3.5	ng/u1	893548	1	7.869	5101570	20.0
1-Hexanol, 2-et...	7.934	28.4	ng/u1	7247110	1	7.869	5101570	20.0
Benzene, 1,2,3-...	7.987	5.7	ng/u1	1444280	1	7.869	5101570	20.0
Oxalic acid, cy...	8.098	12.7	ng/u1	3240390	1	7.869	5101570	20.0
(DEL) Alkane: C...	8.169	5.0	ng/u1	1275220	1	7.869	5101570	20.0
Cyclohexanone, ...	8.298	7.8	ng/u1	1988700	1	7.869	5101570	20.0
Benzene, 1-meth...	8.416	10.8	ng/u1	2743580	1	7.869	5101570	20.0
(DEL) Alkane: S...	8.475	4.8	ng/u1	1236170	1	7.869	5101570	20.0
(DEL) Alkane: S...	8.651	10.0	ng/u1	2560570	1	7.869	5101570	20.0
Benzene, 2-ethy...	8.828	4.2	ng/u1	1080270	1	7.869	5101570	20.0
o-Cymene	8.875	5.3	ng/u1	1355430	1	7.869	5101570	20.0
Benzene, 4-ethy...	8.981	12.3	ng/u1	3141800	1	7.869	5101570	20.0
(DEL) Alkane: S...	9.110	44.2	ng/u1	11281200	1	7.869	5101570	20.0
2-Butanone, 4-p...	9.234	9.6	ng/u1	2444610	1	7.869	5101570	20.0
(DEL) Alkane: C...	9.675	2.0	ng/u1	1340540	2	10.663	13091100	20.0
(DEL) Alkane: S...	9.816	2.9	ng/u1	1924970	2	10.663	13091100	20.0
(DEL) Alkane: S...	9.963	3.1	ng/u1	2037370	2	10.663	13091100	20.0
2,4-Dipropyl-5-...	11.057	8.9	ng/u1	5811820	2	10.663	13091100	20.0
Oxalic acid, 2-...	11.175	4.2	ng/u1	2762490	2	10.663	13091100	20.0
(DEL) Alkane: S...	11.598	2.5	ng/u1	1648830	2	10.663	13091100	20.0
unknown-02	11.939	5.1	ng/u1	3318290	2	10.663	13091100	20.0
(DEL) Alkane: S...	12.104	5.9	ng/u1	3884420	2	10.663	13091100	20.0
(DEL) Alkane: S...	12.916	3.8	ng/u1	1246750	3	14.492	6563240	20.0
(DEL) Alkane: S...	13.057	3.9	ng/u1	1290770	3	14.492	6563240	20.0
(DEL) Alkane: S...	13.339	14.1	ng/u1	4628380	3	14.492	6563240	20.0
Naphthalene, 2,...	13.663	4.0	ng/u1	1320590	3	14.492	6563240	20.0
Oxalic acid, 2-...	13.998	7.3	ng/u1	2389860	3	14.492	6563240	20.0
(DEL) Alkane: S...	14.410	14.5	ng/u1	4752840	3	14.492	6563240	20.0
unknown-03	15.033	4.7	ng/u1	1529330	3	14.492	6563240	20.0
(DEL) Alkane: S...	15.363	12.3	ng/u1	4033400	3	14.492	6563240	20.0
(DEL) Alkane: S...	15.769	6.4	ng/u1	2102550	3	14.492	6563240	20.0
(DEL) Alkane: S...	16.221	16.0	ng/u1	4944540	4	17.233	6169710	20.0
Sulfurous acid,...	16.245	5.1	ng/u1	1572590	4	17.233	6169710	20.0
(DEL) Alkane: S...	17.010	10.1	ng/u1	3114290	4	17.233	6169710	20.0

(DEL) Alkane: S...	17.063	7.5	ng/ul	2308650	4	17.233	6169710	20.0
(DEL) Alkane: S...	17.745	9.6	ng/ul	2946920	4	17.233	6169710	20.0
Pentadecanoic a...	17.915	5.3	ng/ul	1634470	4	17.233	6169710	20.0
(DEL) Alkane: S...	18.433	7.4	ng/ul	2282450	4	17.233	6169710	20.0
(DEL) Alkane: S...	19.068	10.7	ng/ul	3300240	4	17.233	6169710	20.0
(DEL) Alkane: S...	19.645	4.1	ng/ul	1264680	5	21.456	6221900	20.0
(DEL) Alkane: S...	20.174	3.5	ng/ul	1104980	5	21.456	6221900	20.0
Carbonic acid, ...	21.156	9.9	ng/ul	3082240	5	21.456	6221900	20.0
(DEL) Alkane: S...	21.680	2.8	ng/ul	879826	5	21.456	6221900	20.0
Dodecyl pentyl ...	22.139	5.1	ng/ul	1575440	5	21.456	6221900	20.0
Oxalic acid, 2-...	22.262	4.2	ng/ul	1296330	5	21.456	6221900	20.0

Instrument :

BNA_M

ClientSampleId :

DCVY6DL2