

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
 Data File : BM047700.D
 Acq On : 28 Sep 2024 20:38
 Operator : RC/JU
 Sample : P3927-08DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDC93DL

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-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047700.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.840	479	485	494	rVB2	2363669	3472822	5.60%	1.130%
2	6.381	573	577	582	rBV	751158	1094161	1.76%	0.356%
3	6.463	582	591	594	rVV3	586919	1348441	2.17%	0.439%
4	6.816	640	651	656	rBV3	1167041	2800950	4.51%	0.911%
5	6.875	656	661	664	rBV	1425150	2083158	3.36%	0.678%
6	6.910	664	667	672	rVB	1561899	2201401	3.55%	0.716%
7	6.981	672	679	685	rBV3	1790379	3762286	6.06%	1.224%
8	7.046	685	690	695	rVB	948795	1400841	2.26%	0.456%
9	7.210	712	718	721	rBV	953823	1465577	2.36%	0.477%
10	7.246	721	724	728	rVV	1169065	1459721	2.35%	0.475%
11	7.451	753	759	767	rVB2	8211792	14002259	22.57%	4.554%
12	7.804	814	819	825	rVB2	2038531	3521700	5.68%	1.145%
13	7.893	825	834	838	rBV3	542878	1592047	2.57%	0.518%
14	7.934	838	841	850	rVV3	895090	1608890	2.59%	0.523%
15	8.016	850	855	857	rVV2	678107	1083198	1.75%	0.352%
16	8.040	857	859	863	rVV	848928	1145925	1.85%	0.373%
17	8.087	863	867	869	rVV	551762	905266	1.46%	0.294%
18	8.116	869	872	878	rVB4	667528	1095231	1.77%	0.356%
19	8.245	889	894	899	rBV	1364707	2308597	3.72%	0.751%
20	8.298	899	903	908	rVV2	700988	1544563	2.49%	0.502%
21	8.357	908	913	919	rVV2	1549665	3046931	4.91%	0.991%
22	8.416	919	923	926	rVV	1091767	1516471	2.44%	0.493%
23	8.457	926	930	932	rVV2	1349207	2013230	3.25%	0.655%
24	8.487	932	935	945	rVB	1887247	3360566	5.42%	1.093%
25	8.593	947	953	956	rBV	1534377	2541780	4.10%	0.827%
26	8.622	956	958	967	rVB	808413	1261076	2.03%	0.410%
27	8.769	979	983	987	rBV	675243	989195	1.59%	0.322%
28	8.816	987	991	1000	rVB2	945670	1774825	2.86%	0.577%
29	8.922	1000	1009	1020	rBV2	1175457	2998201	4.83%	0.975%
30	9.051	1020	1031	1037	rBV	7170750	11249477	18.13%	3.659%
31	9.175	1042	1052	1058	rBV8	754504	2133137	3.44%	0.694%
32	9.251	1058	1065	1069	rBV4	616932	1331693	2.15%	0.433%
33	9.616	1119	1127	1133	rVB2	487799	1076509	1.74%	0.350%
34	9.704	1133	1142	1146	rBV6	432861	1166337	1.88%	0.379%
35	9.757	1146	1151	1155	rBV3	650663	1303066	2.10%	0.424%
36	9.898	1166	1175	1180	rVB	806057	1907733	3.07%	0.621%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	9.969	1181	1187	1191	rBV	610538	970296	1.56%	0.316%
38	10.045	1197	1200	1204	rVV	697258	1140773	1.84%	0.371%
39	10.151	1213	1218	1223	rBV2	525863	913738	1.47%	0.297%
40	10.310	1239	1245	1250	rBV2	637551	1037897	1.67%	0.338%

41	10.604	1286	1295	1301	rVV3	8318638	15720918	25.34%	5.113%
42	10.663	1301	1305	1311	rVV	602468	1261520	2.03%	0.410%
43	10.734	1311	1317	1323	rVV3	2377604	4350601	7.01%	1.415%
44	10.992	1355	1361	1369	rBV	2409062	3999721	6.45%	1.301%
45	11.122	1375	1383	1389	rVB	831359	1672731	2.70%	0.544%

46	11.504	1443	1448	1450	rVV2	503099	938194	1.51%	0.305%
47	11.534	1451	1453	1467	rVV5	585035	1130019	1.82%	0.368%
48	11.645	1467	1472	1476	rVV	1224312	2000272	3.22%	0.651%
49	11.710	1476	1483	1490	rVB	1763560	3276395	5.28%	1.066%
50	11.881	1503	1512	1520	rBV2	1705265	2976623	4.80%	0.968%

51	12.039	1533	1539	1542	rBV	2254820	3577932	5.77%	1.164%
52	12.075	1542	1545	1550	rVB	1549759	2185868	3.52%	0.711%
53	12.286	1573	1581	1590	rVB2	2376381	4852467	7.82%	1.578%
54	12.445	1603	1608	1612	rBV3	1098837	1838111	2.96%	0.598%
55	12.492	1612	1616	1620	rVB	576363	874269	1.41%	0.284%

56	12.686	1643	1649	1661	rBV4	735177	1800051	2.90%	0.585%
57	12.857	1673	1678	1687	rVB3	383573	858165	1.38%	0.279%
58	13.286	1745	1751	1757	rVV	2526828	3691296	5.95%	1.201%
59	13.504	1781	1788	1796	rVB4	575002	1362300	2.20%	0.443%
60	13.633	1801	1810	1815	rBV5	716770	1992371	3.21%	0.648%

61	13.716	1820	1824	1828	rBV	734094	1015038	1.64%	0.330%
62	13.769	1828	1833	1837	rVV	784226	1549115	2.50%	0.504%
63	13.822	1837	1842	1845	rVV3	800256	1670302	2.69%	0.543%
64	13.945	1860	1863	1867	rVV2	784729	1040835	1.68%	0.339%
65	14.086	1883	1887	1898	rVB3	1072559	1761474	2.84%	0.573%

66	14.363	1928	1934	1938	rBV	9734680	13171612	21.23%	4.284%
67	14.445	1943	1948	1954	rVV	1613552	2525643	4.07%	0.821%
68	14.527	1958	1962	1968	rVV5	674879	1101763	1.78%	0.358%
69	14.592	1968	1973	1980	rVV	2407798	3242948	5.23%	1.055%
70	14.663	1980	1985	1997	rVB7	492519	1168717	1.88%	0.380%

71	14.845	2013	2016	2022	rVB4	588797	982753	1.58%	0.320%
72	14.916	2022	2028	2032	rBV3	468785	878188	1.42%	0.286%
73	14.980	2034	2039	2042	rBV3	577013	942715	1.52%	0.307%
74	15.069	2049	2054	2059	rVV	5086313	8021009	12.93%	2.609%
75	15.139	2063	2066	2070	rVB	759809	915895	1.48%	0.298%

76	15.210	2073	2078	2086	rVV	2640797	4118555	6.64%	1.340%
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Integration Parameters: LSCINT.P

Integrator: RTE

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77	15.304	2086	2094	2099	rVB	2650958	4412062	7.11%	1.435%
78	15.545	2129	2135	2140	rBV	889159	1433568	2.31%	0.466%
79	15.710	2158	2163	2169	rVB	843429	1397630	2.25%	0.455%
80	15.804	2175	2179	2185	rVB4	711294	1074713	1.73%	0.350%
81	15.921	2195	2199	2209	rVB5	372802	884898	1.43%	0.288%
82	16.080	2222	2226	2230	rVB2	832069	1131740	1.82%	0.368%
83	16.169	2236	2241	2244	rBV	2678453	3938209	6.35%	1.281%
84	16.857	2348	2358	2367	rBV	28505288	62040727	100.00%	20.179%
85	16.957	2371	2375	2379	rVV	2204656	2809856	4.53%	0.914%
86	17.010	2379	2384	2393	rVB	1095359	1861095	3.00%	0.605%
87	17.186	2410	2414	2419	rBV	1877194	2762836	4.45%	0.899%
88	17.286	2427	2431	2435	rVB	635366	899298	1.45%	0.293%
89	17.480	2460	2464	2472	rVB4	595034	1158960	1.87%	0.377%
90	17.692	2495	2500	2505	rVB	2023623	2463817	3.97%	0.801%
91	17.863	2524	2529	2533	rVB	1897124	2230383	3.60%	0.725%
92	18.380	2613	2617	2623	rBV	1450302	1821599	2.94%	0.592%
93	19.027	2720	2727	2731	rVB2	1631053	3216840	5.19%	1.046%
94	19.162	2746	2750	2758	rVB	742755	966321	1.56%	0.314%
95	19.586	2820	2822	2834	rVB	918671	1172169	1.89%	0.381%
96	20.121	2909	2913	2922	rBV2	804508	1150857	1.86%	0.374%
97	21.098	3074	3079	3089	rVB	1303385	1823574	2.94%	0.593%
98	21.403	3126	3131	3138	rBV2	1838700	2827292	4.56%	0.920%
99	22.174	3257	3262	3272	rVB2	1049272	1920894	3.10%	0.625%
100	24.409	3635	3642	3661	rVB	1286717	3951887	6.37%	1.285%

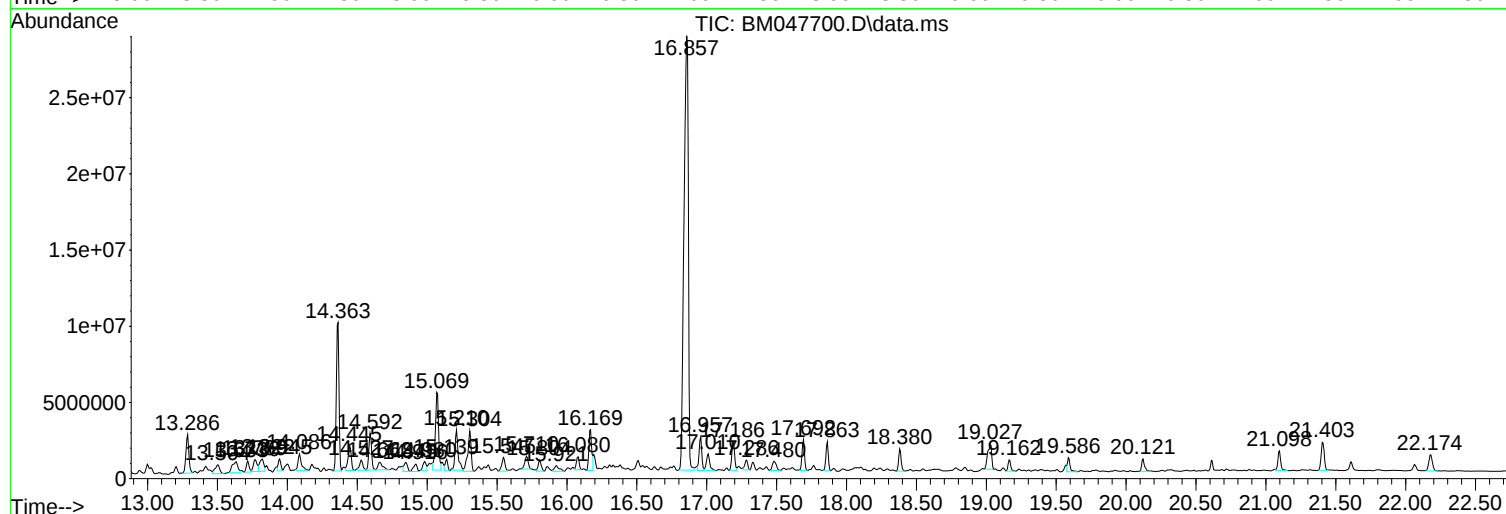
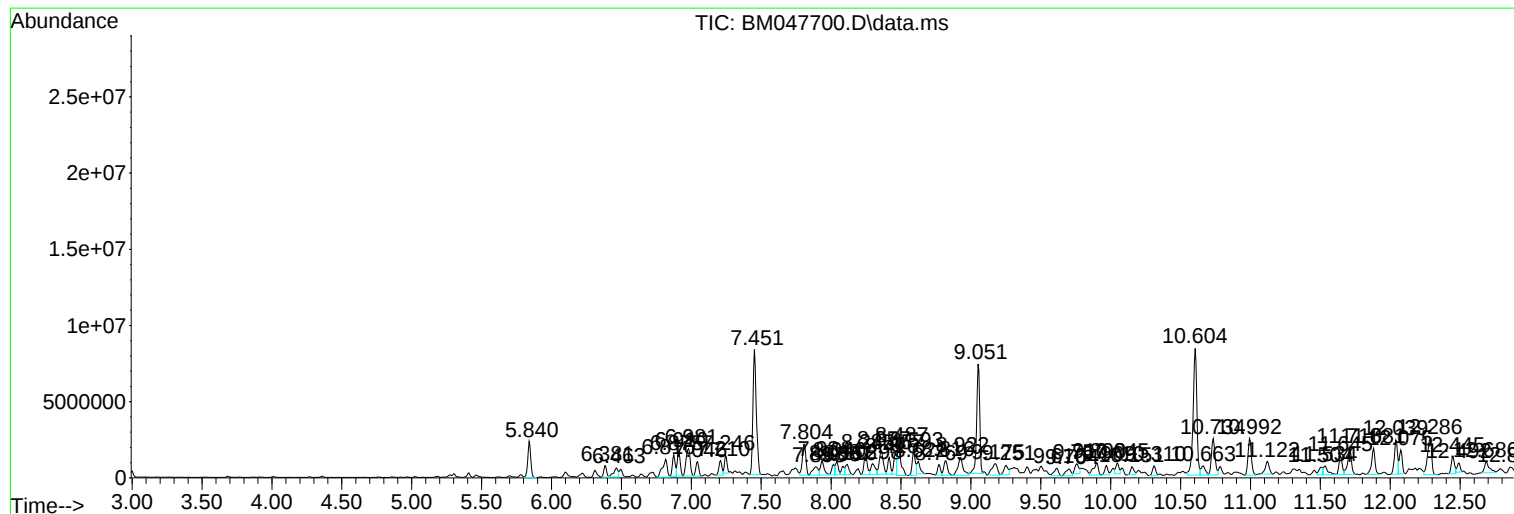
Sum of corrected areas: 307447576

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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

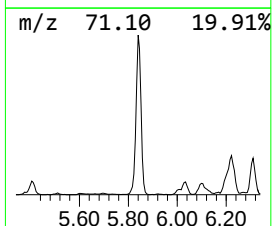
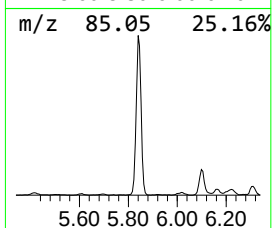
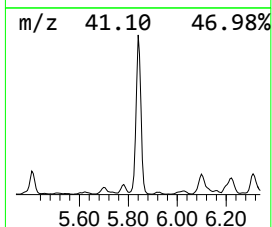
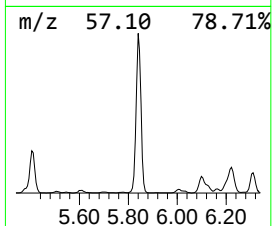
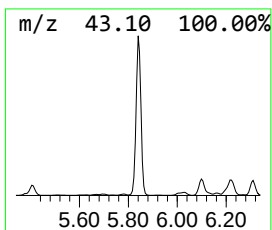
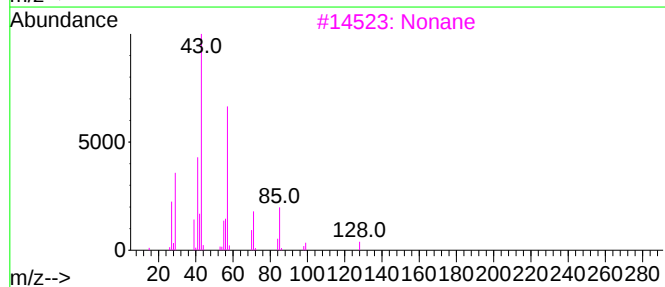
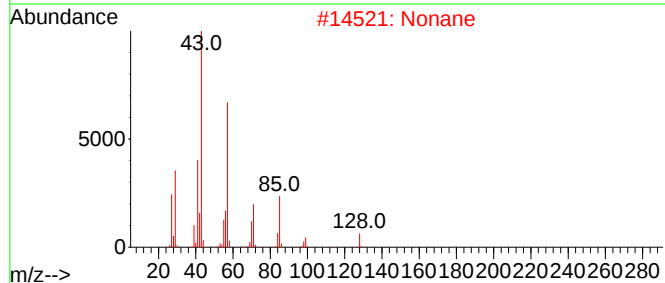
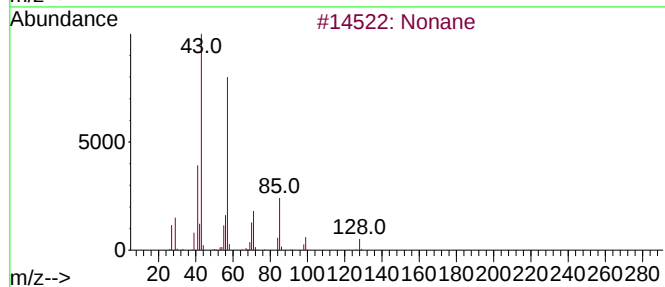
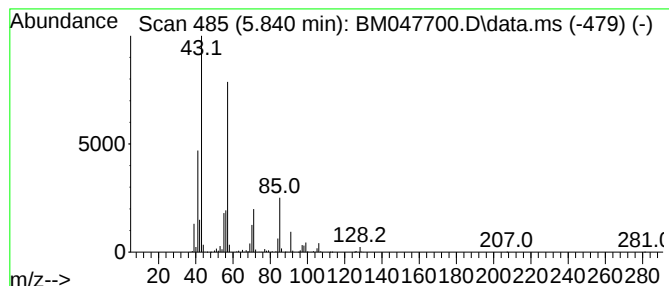
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.840	19.72 ng/ul	3472820	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	87
2	Nonane			128	C9H20	000111-84-2	80
3	Nonane			128	C9H20	000111-84-2	72
4	Nonane			128	C9H20	000111-84-2	72
5	Octane			114	C8H18	000111-65-9	59



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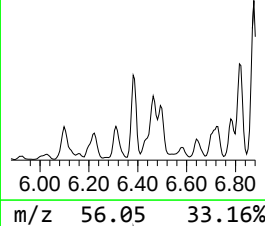
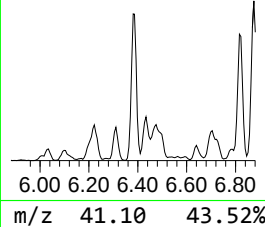
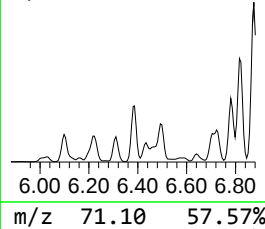
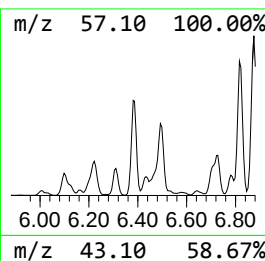
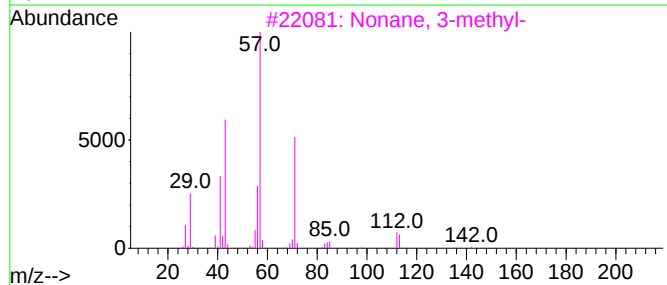
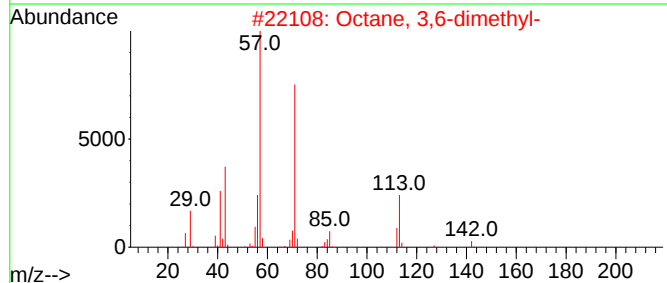
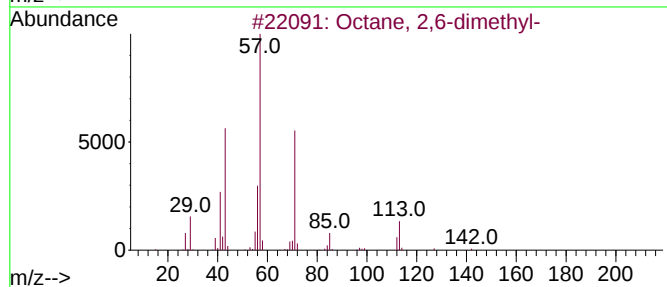
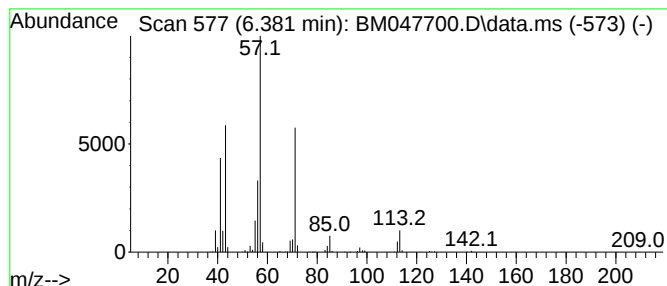
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.381	6.21 ng/ul	1094160	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	90
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
4			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	64
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	58



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Instrument :
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ClientSampleId :
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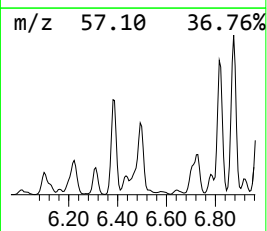
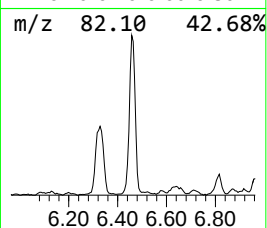
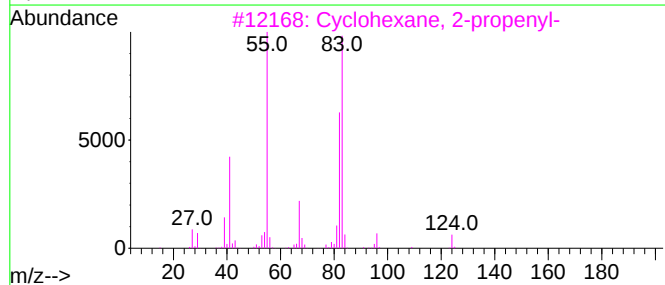
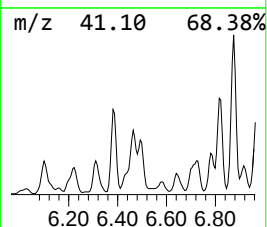
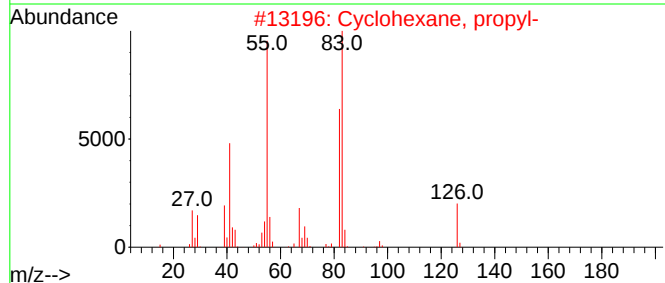
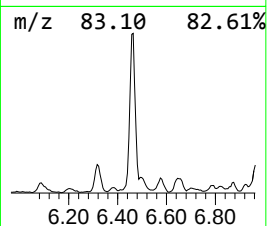
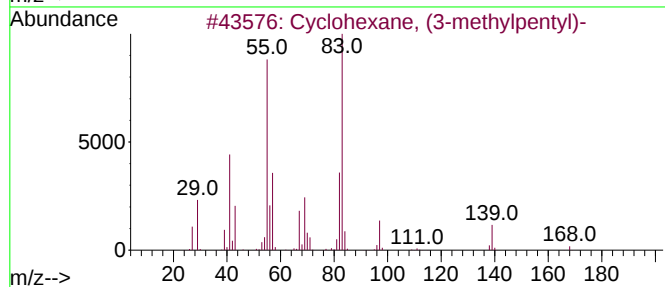
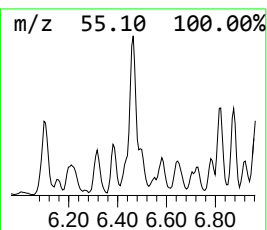
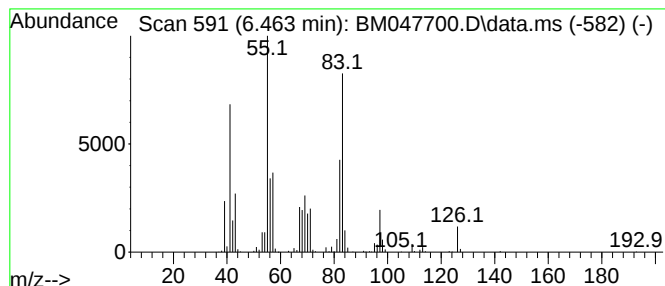
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic6.463 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.463	7.66 ng/ul	1348440	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (3-methylpentyl)-	168	C12H24	061142-38-9	59
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	53
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	50
4			Cyclohexane, undecyl-	238	C17H34	054105-66-7	43
5			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	38



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Data File : BM047700.D
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Operator : RC/JU
Sample : P3927-08DL 5X
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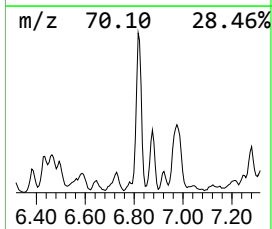
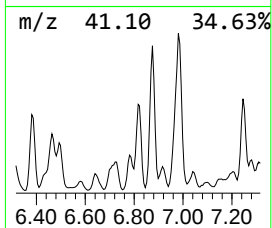
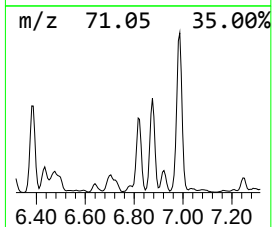
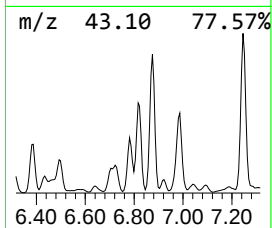
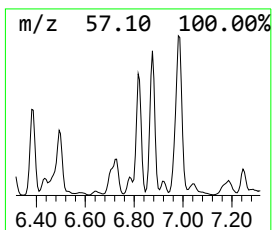
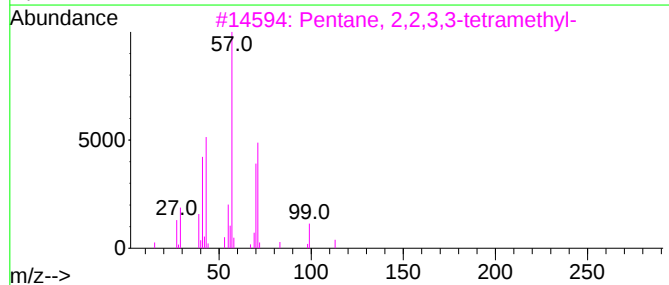
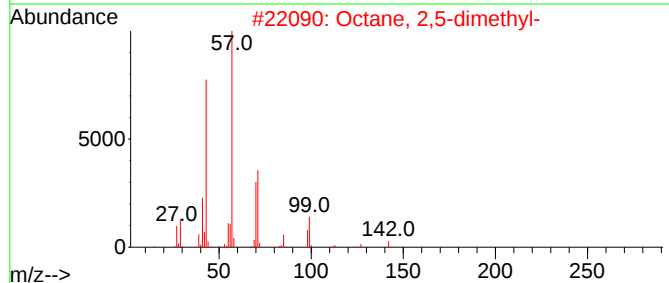
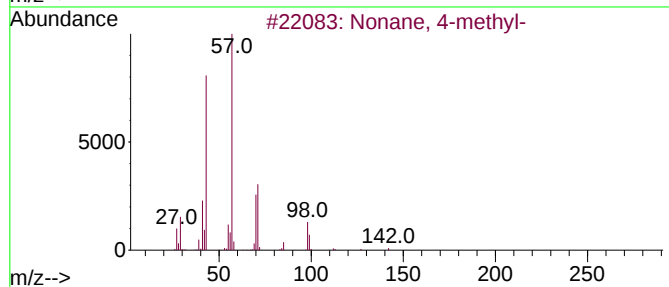
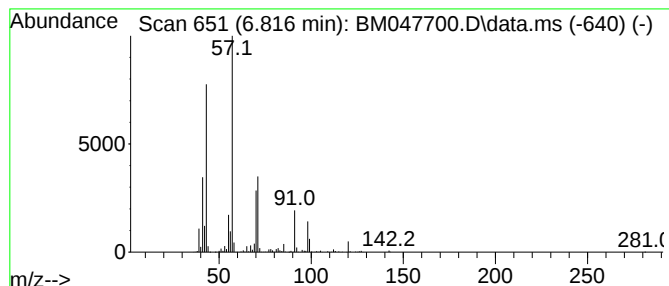
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.816	15.91 ng/ul	2800950	1,4-Dichlorobenzene-d4			7.816
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-		142	C10H22	017301-94-9	91
2	Octane, 2,5-dimethyl-		142	C10H22	015869-89-3	72
3	Pentane, 2,2,3,3-tetramethyl-		128	C9H20	007154-79-2	64
4	Heptane, 2,3,4-trimethyl-		142	C10H22	052896-95-4	64
5	Dodecane, 6-methyl-		184	C13H28	006044-71-9	59



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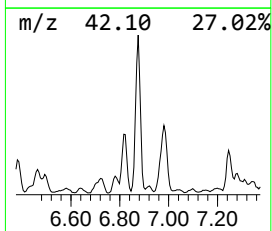
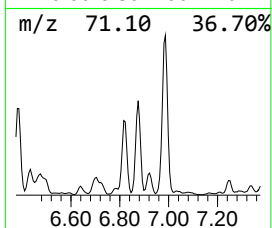
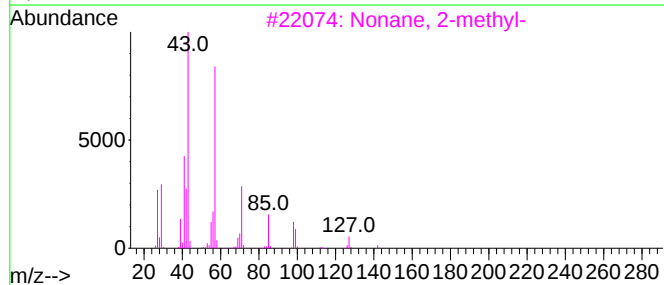
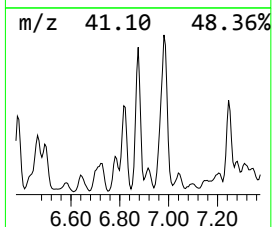
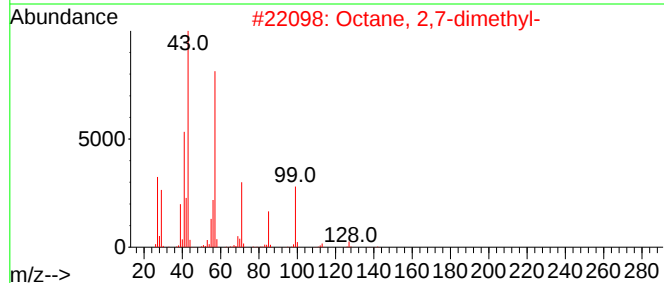
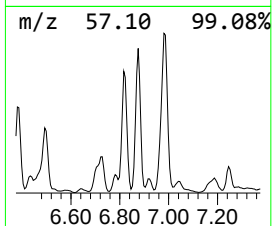
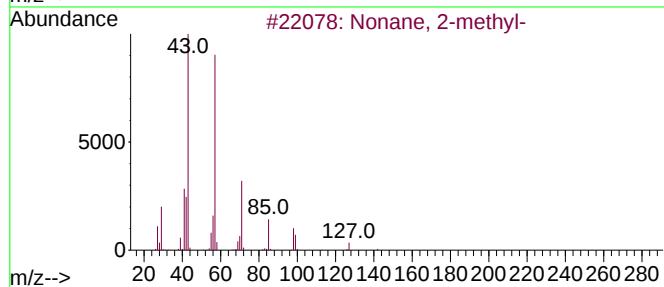
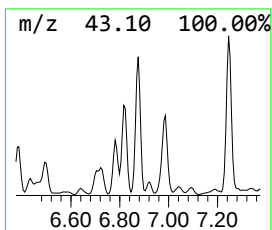
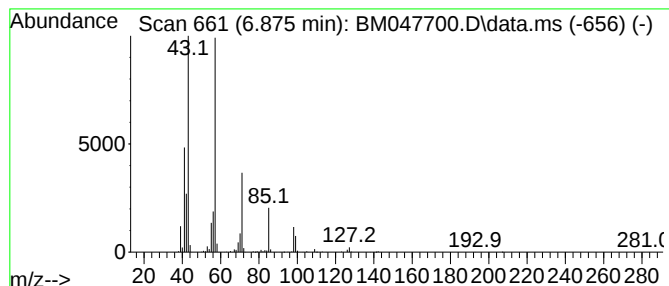
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.875	11.83 ng/ul	2083160	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	91
2			Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	83
3			Nonane, 2-methyl-	142	C10H22	000871-83-0	78
4			Decane	142	C10H22	000124-18-5	64
5			Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	59



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Data File : BM047700.D
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Misc :
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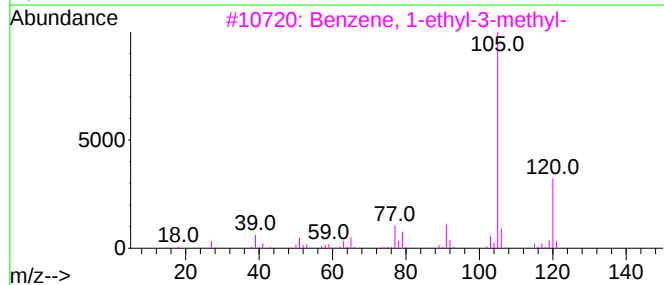
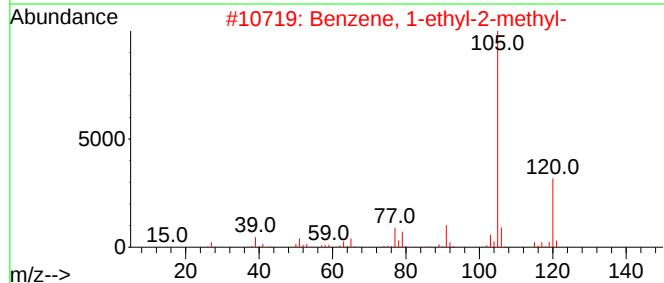
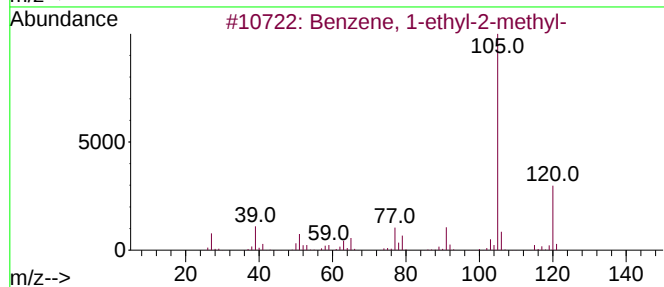
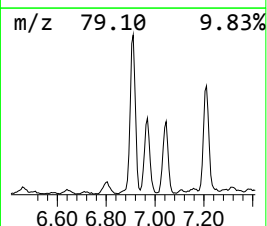
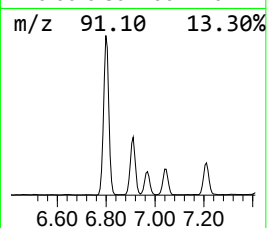
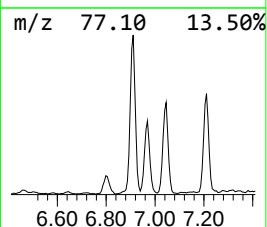
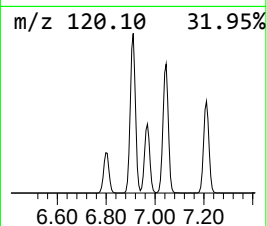
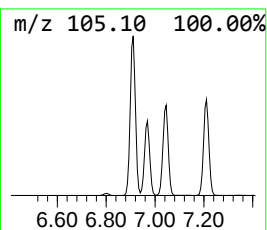
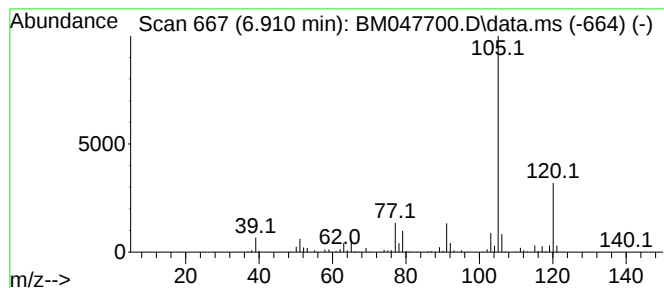
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.910	12.50 ng/ul	2201400	1,4-Dichlorobenzene-d4	7.816

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-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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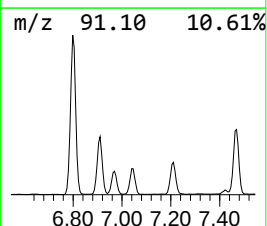
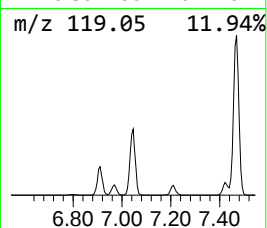
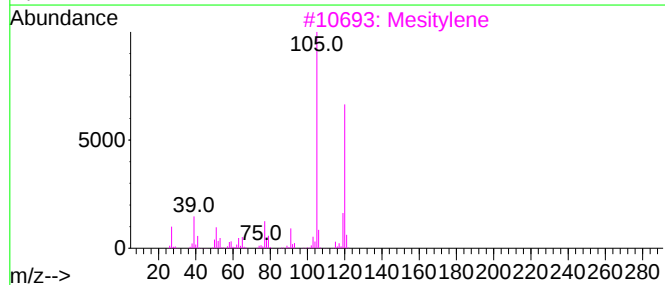
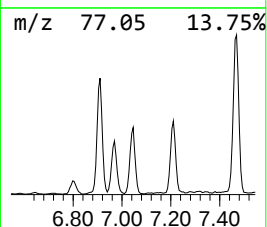
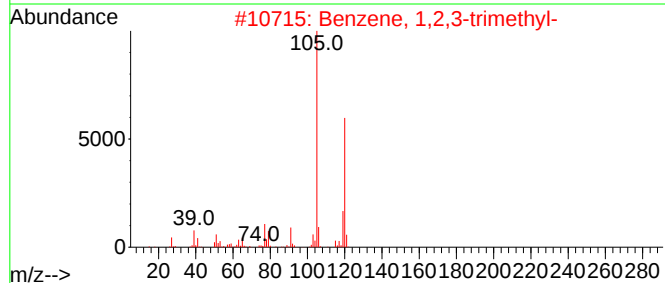
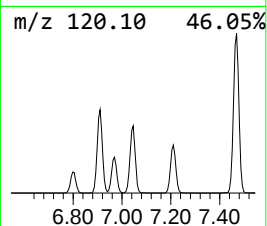
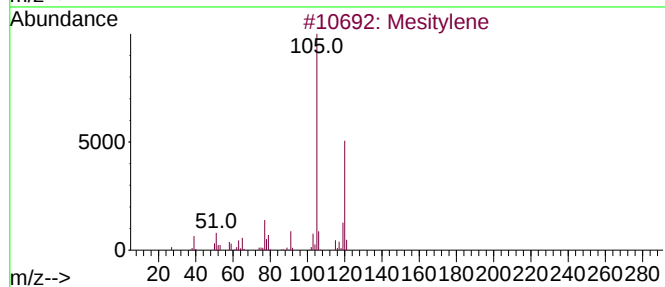
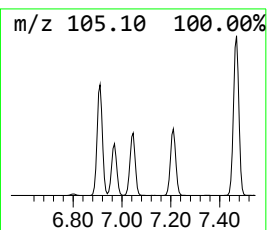
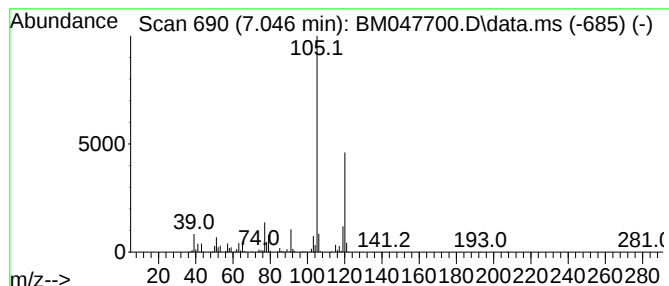
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Mesitylene Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.046	7.96 ng/ul	1400840	1,4-Dichlorobenzene-d4	7.816

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			Mesitylene	120	C9H12	000108-67-8	95
4			Mesitylene	120	C9H12	000108-67-8	95
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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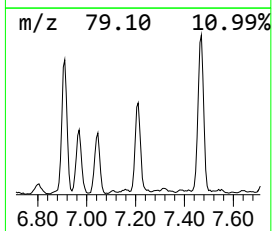
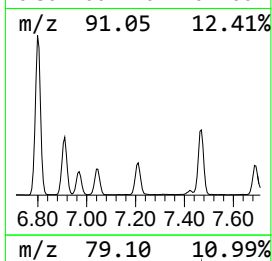
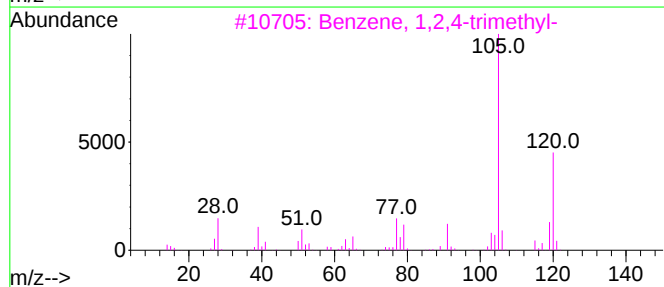
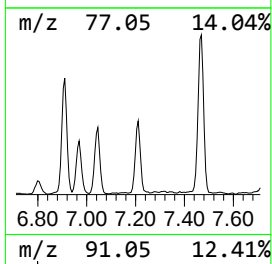
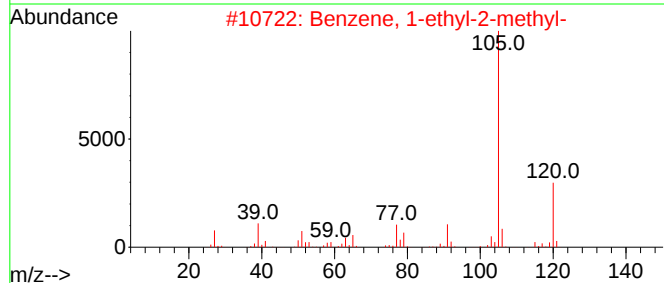
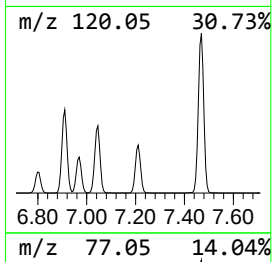
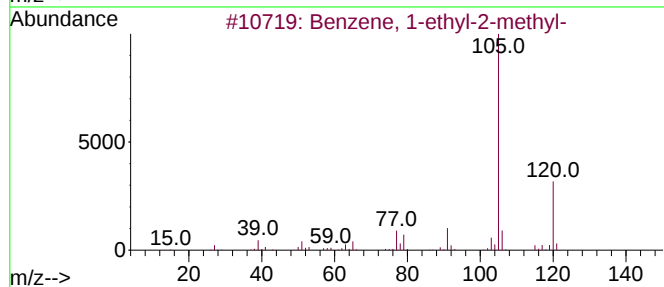
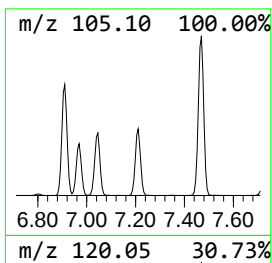
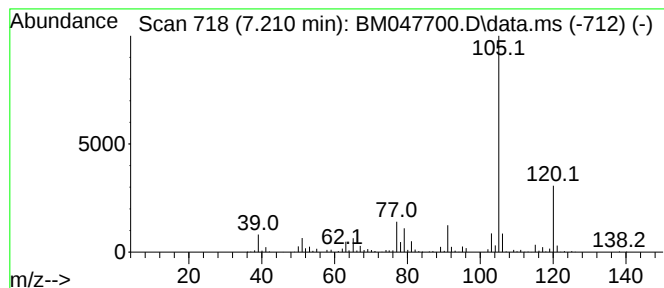
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,4-trimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.210	8.32 ng/ul	1465580	1,4-Dichlorobenzene-d4	7.816

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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
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Misc :
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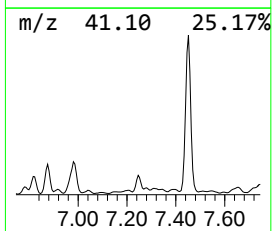
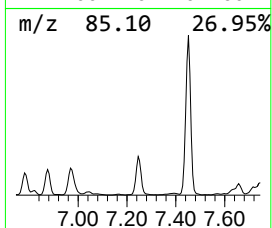
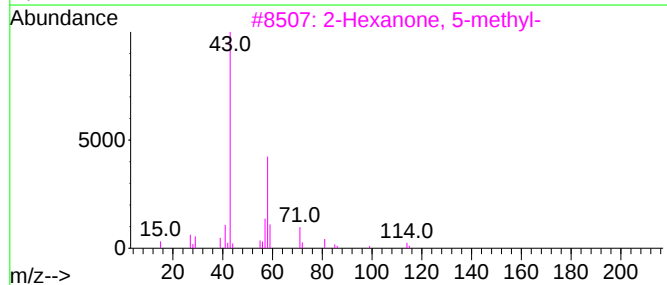
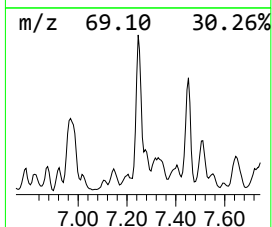
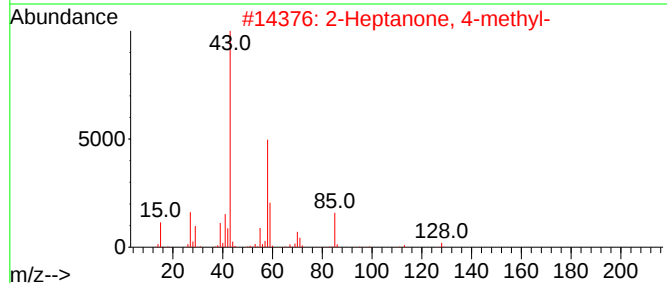
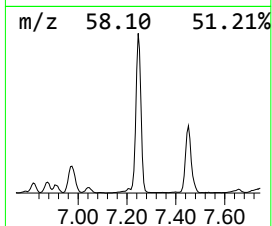
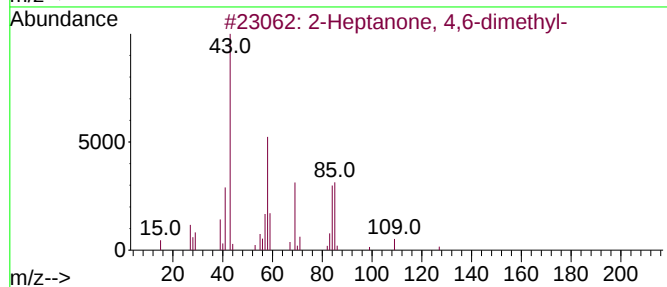
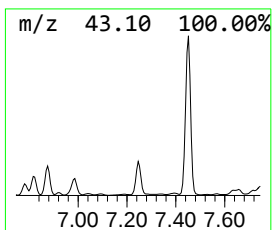
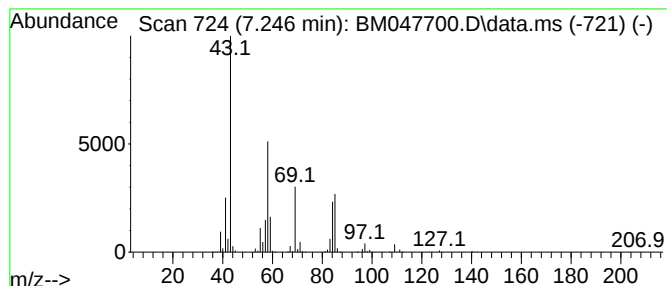
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Heptanone, 4,6-dimethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.246	8.29 ng/ul	1459720	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	90		
2	2-Heptanone, 4-methyl-	128	C8H16O	006137-06-0	38		
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	38		
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	38		
5	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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ALS Vial : 15 Sample Multiplier: 1

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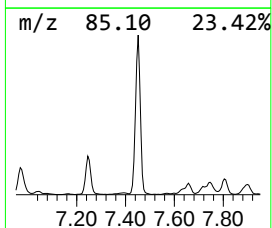
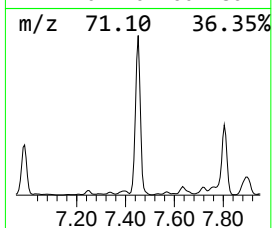
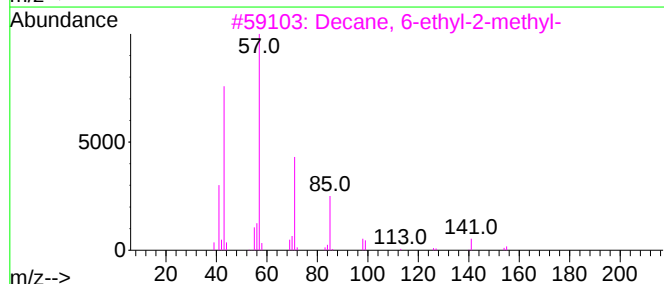
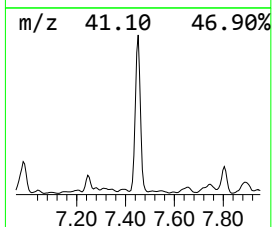
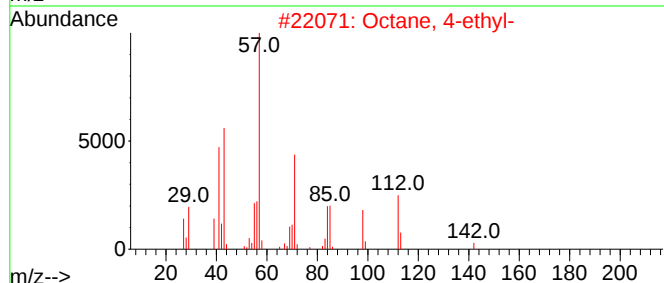
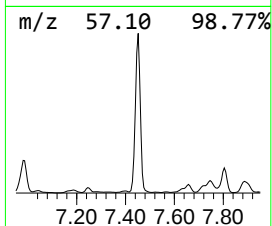
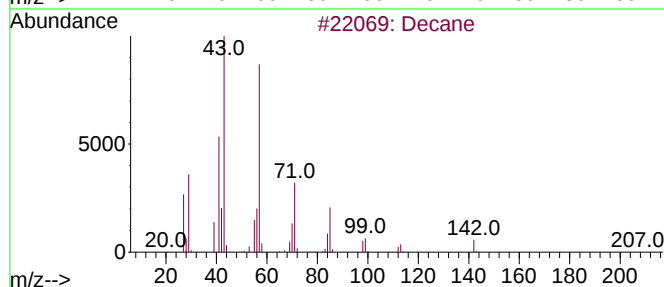
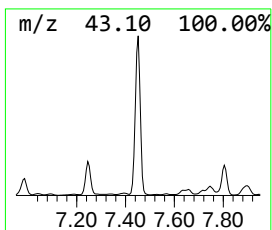
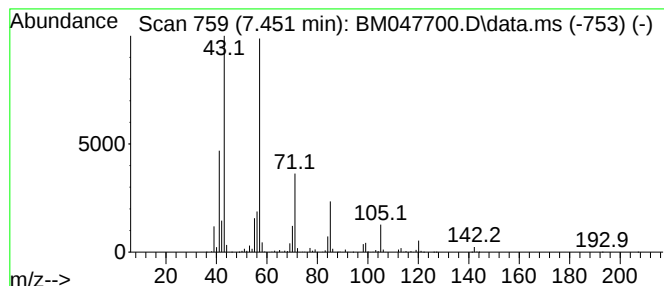
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.451	79.52 ng/ul	14002300	1,4-Dichlorobenzene-d4	7.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane	142	C10H22	000124-18-5	72
2		Octane, 4-ethyl-	142	C10H22	015869-86-0	64
3		Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64
4		Hexatriacontane	507	C36H74	000630-06-8	64
5		Undecane	156	C11H24	001120-21-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

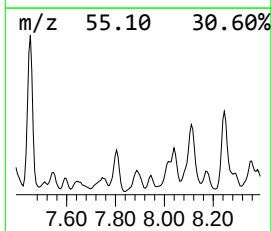
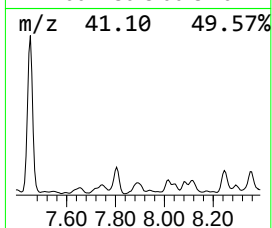
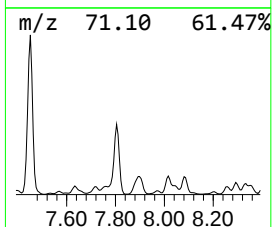
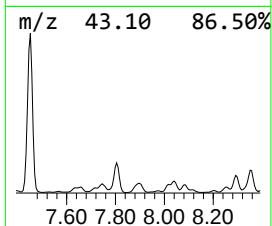
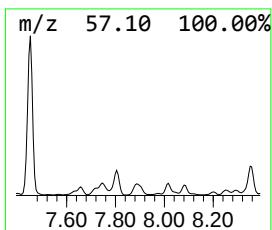
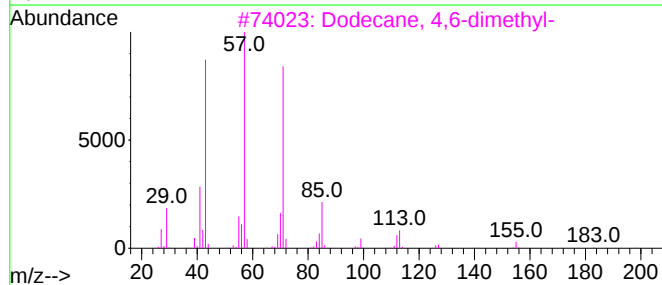
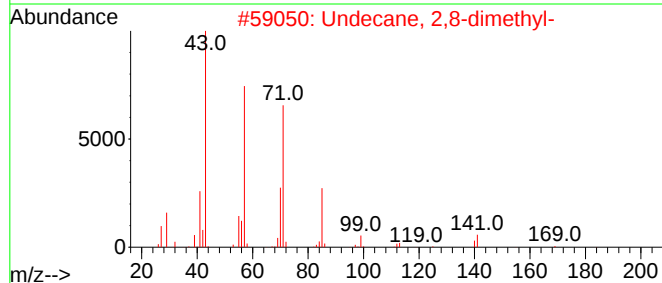
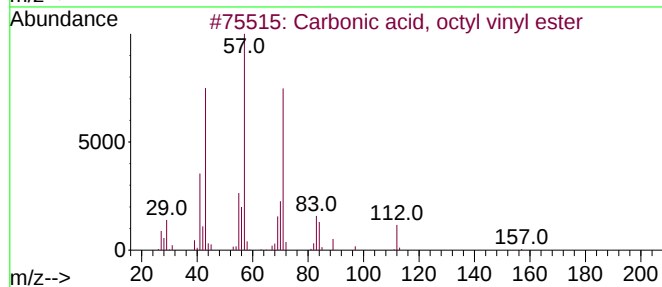
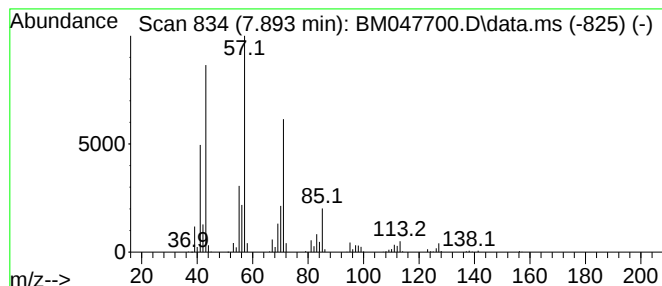
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Carbonic acid, octyl vinyl ... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.893	9.04 ng/ul	1592050	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	59
2			Undecane, 2,8-dimethyl-	184	C13H28	017301-25-6	59
3			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	53
4			Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	53
5			Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

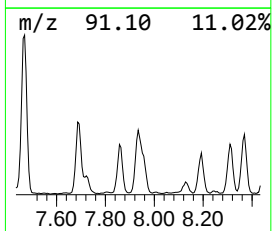
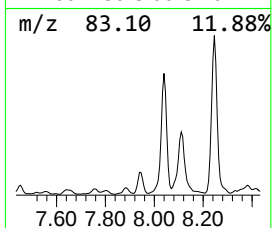
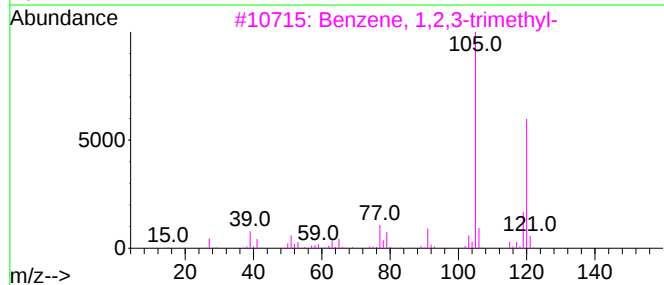
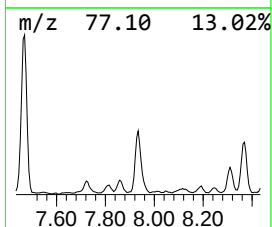
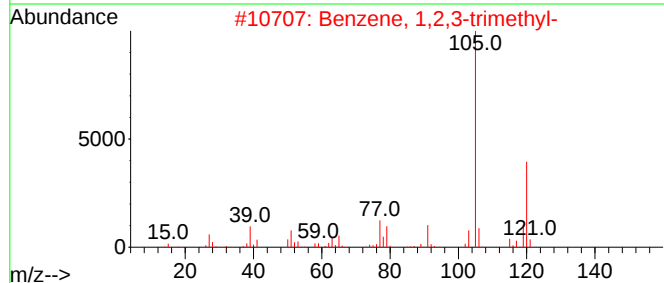
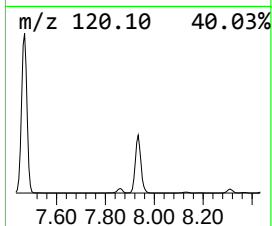
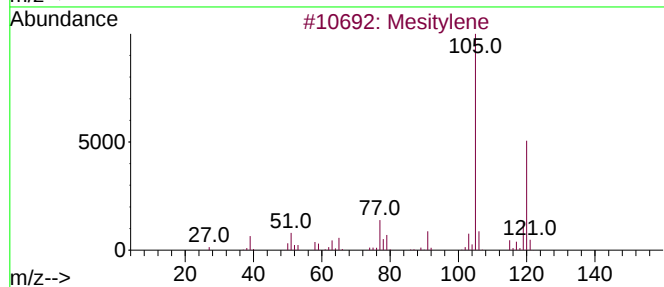
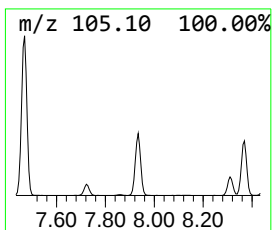
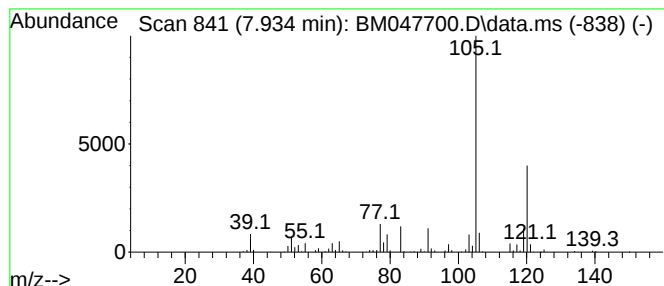
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	9.14 ng/ul	1608890	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	95
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	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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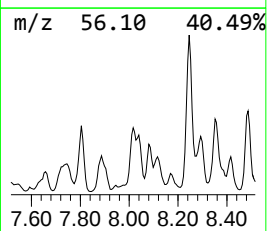
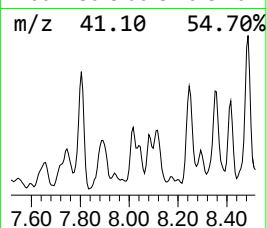
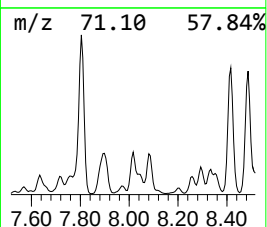
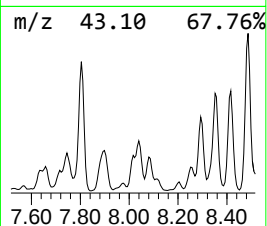
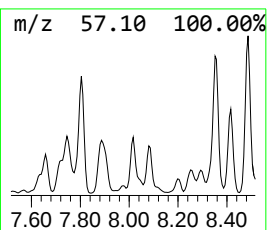
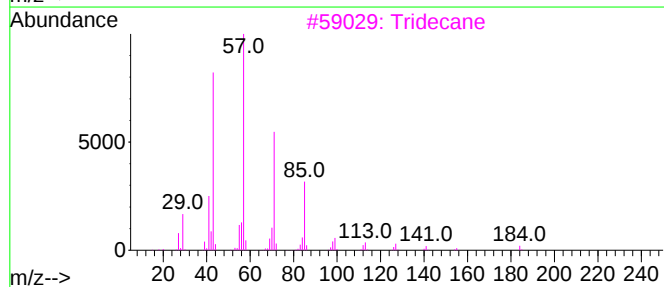
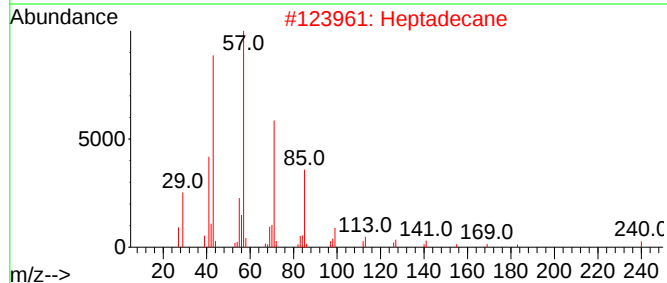
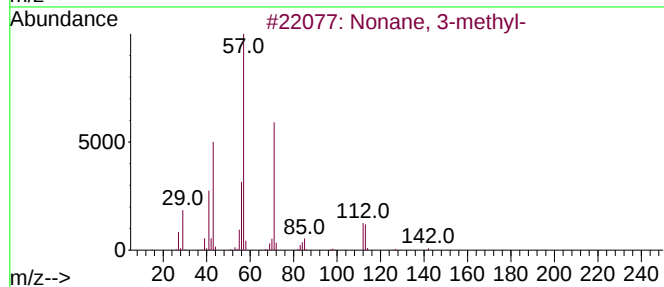
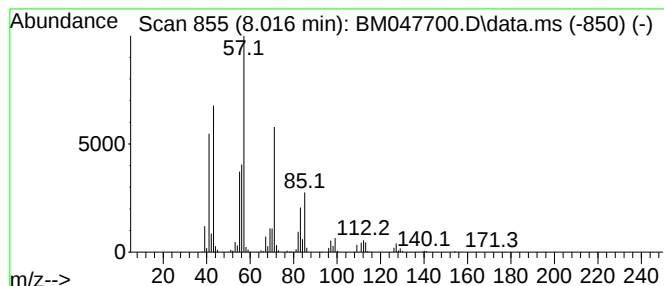
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.016	6.15 ng/ul	1083200	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	53
2			Heptadecane	240	C17H36	000629-78-7	49
3			Tridecane	184	C13H28	000629-50-5	49
4			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	49
5			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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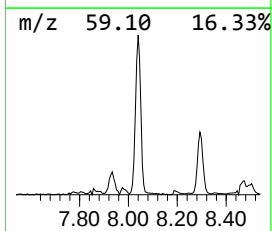
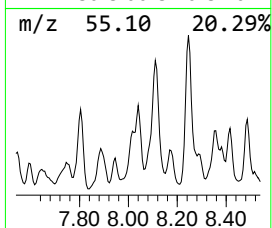
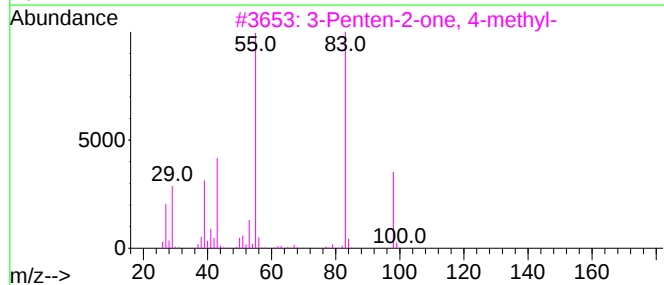
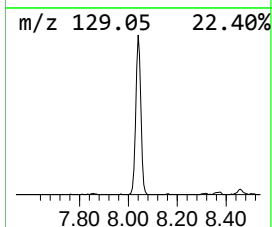
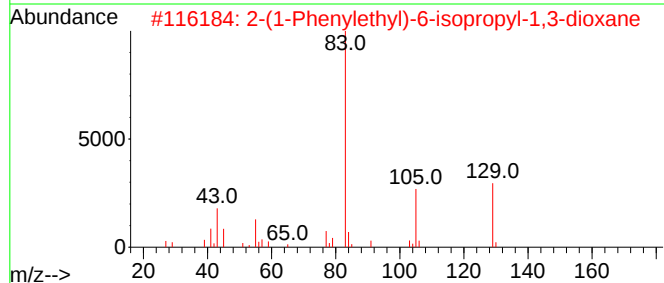
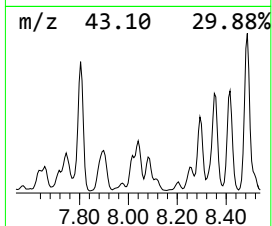
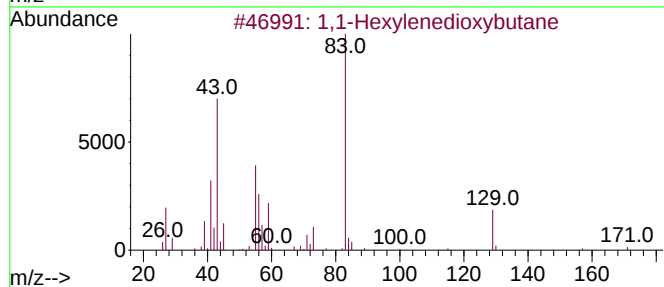
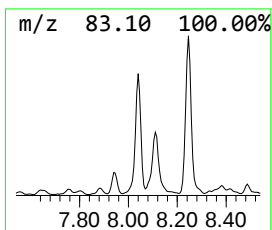
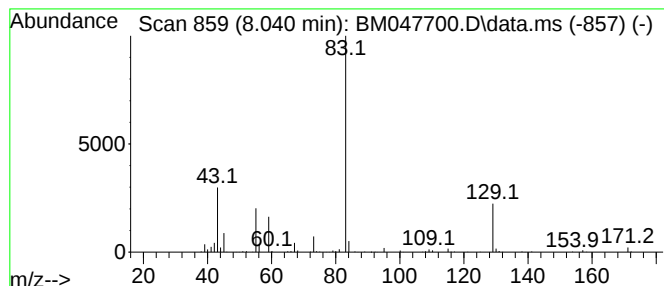
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-01 Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.040	6.51 ng/ul	1145930	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	40
2			2-(1-Phenylethyl)-6-isopropyl-1,...	234	C15H22O2	1000431-40-0	39
3			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	9
4			1,2-Benzenediol, o-(3-methylbut-...	192	C11H12O3	1000325-91-3	9
5			Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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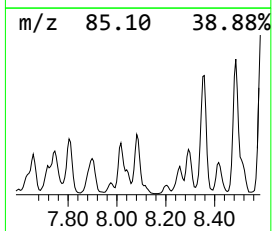
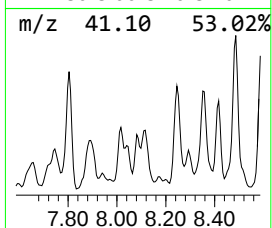
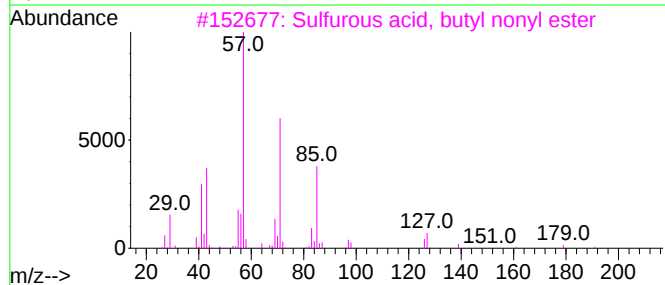
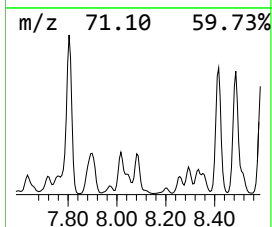
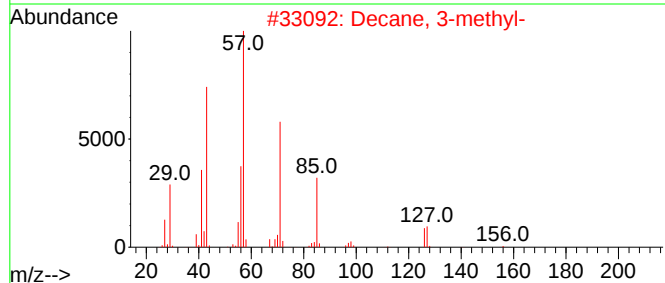
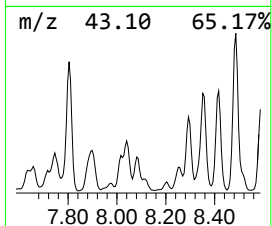
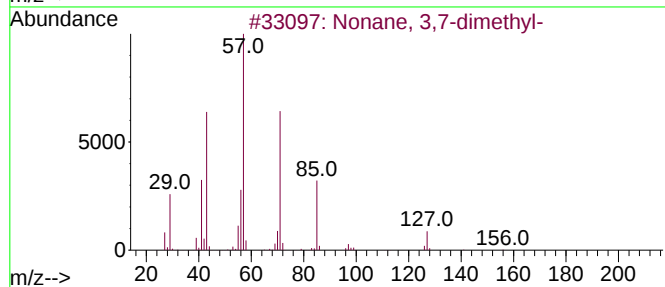
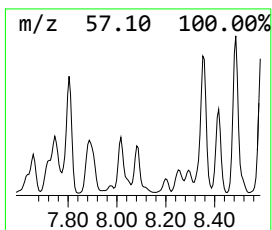
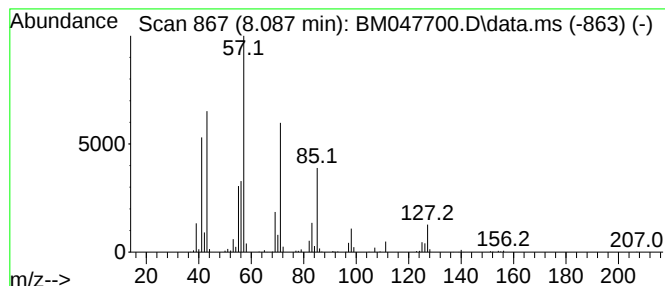
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.087	5.14 ng/ul	905266	1,4-Dichlorobenzene-d4	7.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	90
2		Decane, 3-methyl-	156	C11H24	013151-34-3	87
3		Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	64
4		Dodecane, 3-methyl-	184	C13H28	017312-57-1	64
5		Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
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Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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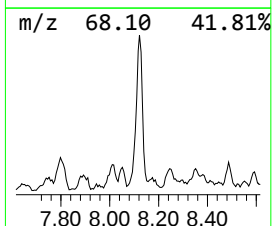
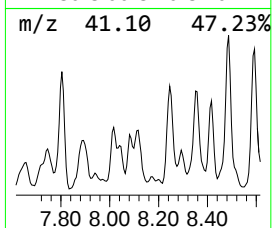
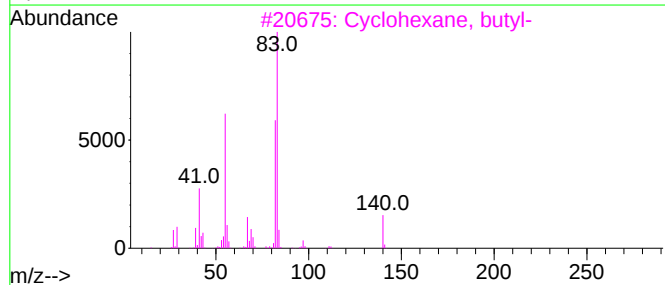
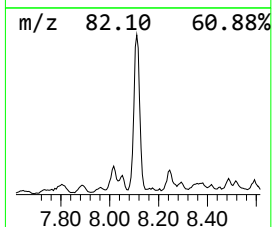
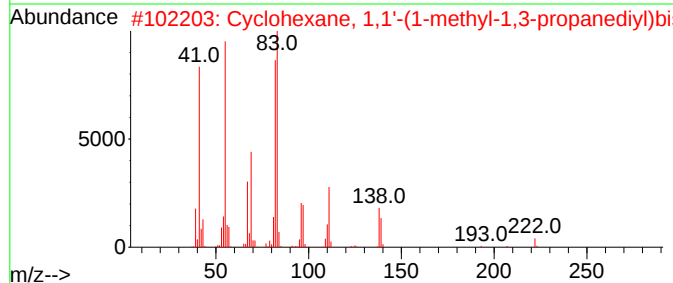
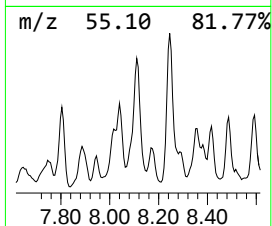
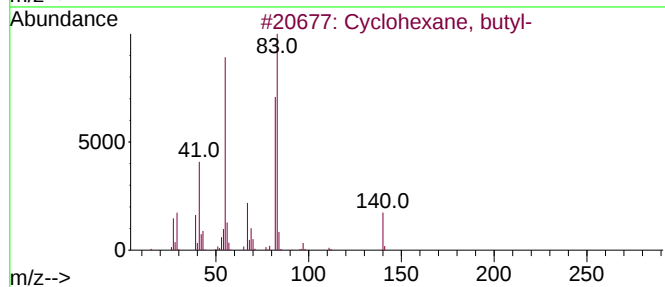
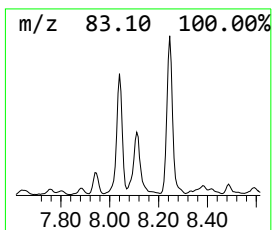
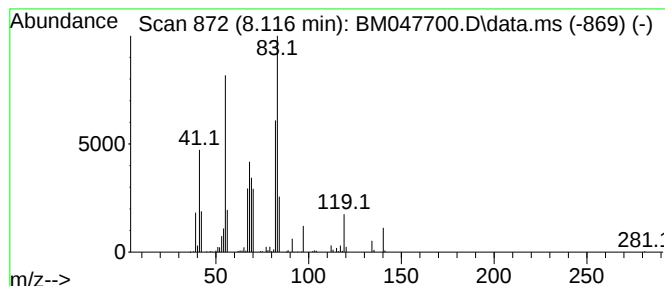
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Cyclic8.116 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.116	6.22 ng/ul	1095230	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	47
2			Cyclohexane, 1,1'-(1-methyl-1,3-...	222	C16H30	041851-35-8	47
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	47
4			Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	45
5			5-Bromo-2-(4-piperidinyloxy)pyri...	256	C10H13BrN2O	194668-50-3	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

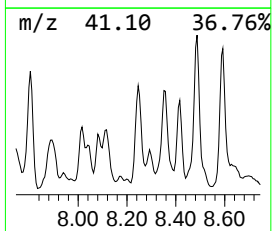
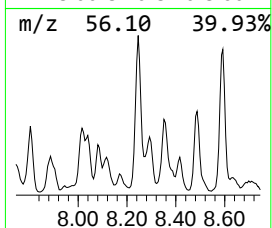
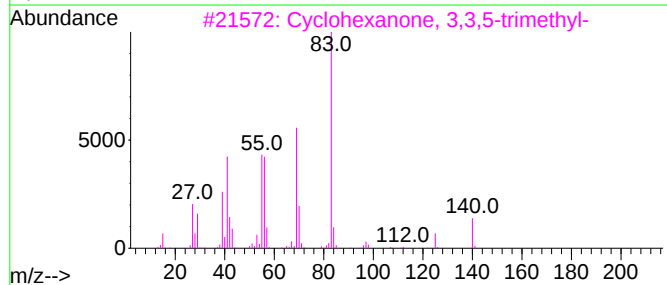
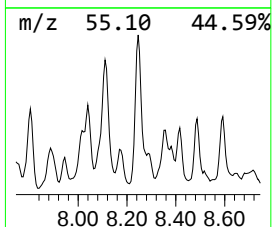
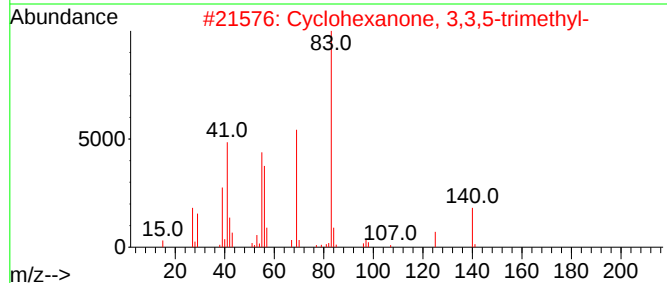
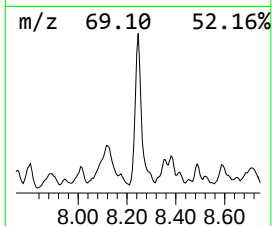
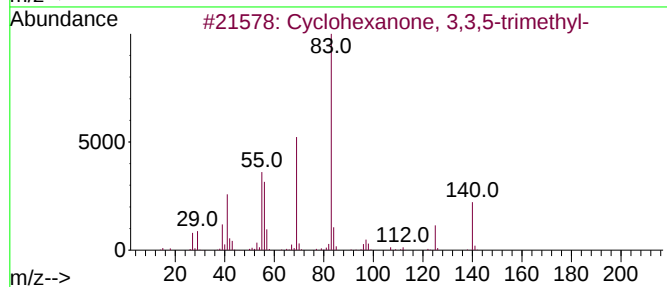
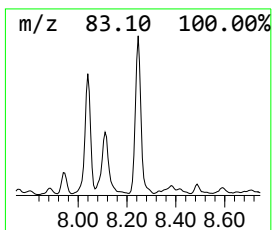
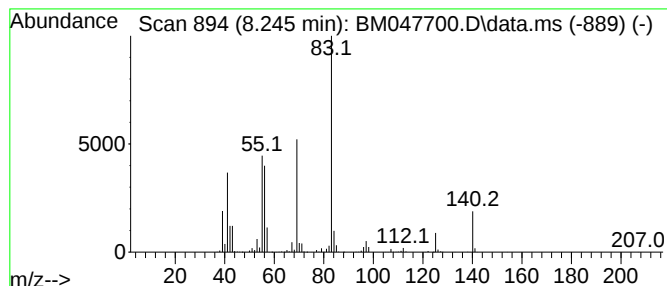
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Cyclohexanone, 3,3,5-trimet... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.245	13.11 ng/ul	2308600	1,4-Dichlorobenzene-d4			7.816
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	94
3	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	91
4	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	90
5	1,4-Pentadien-3-ol		84	C5H8O	000922-65-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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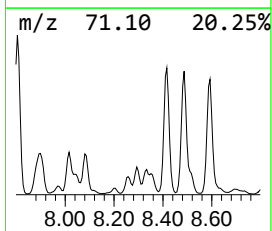
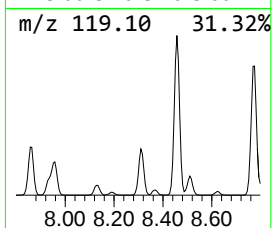
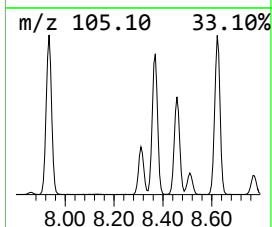
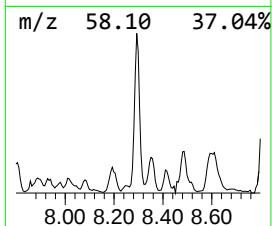
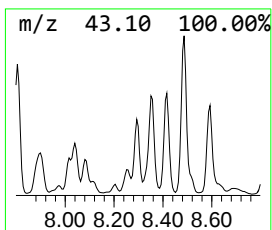
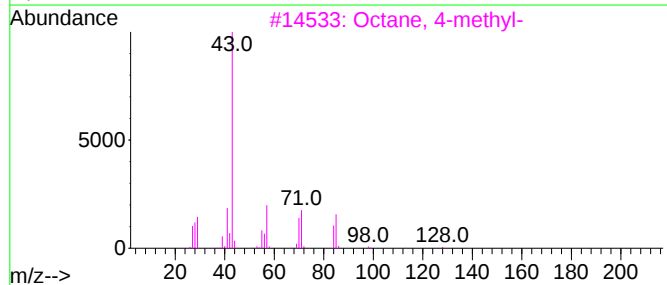
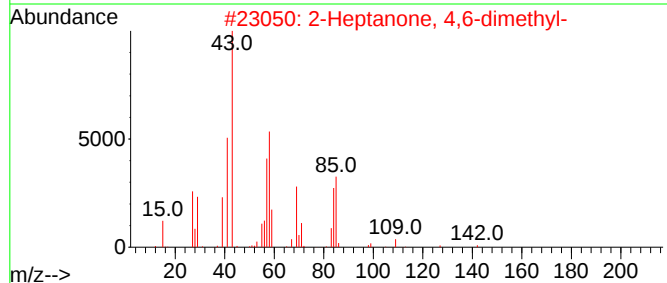
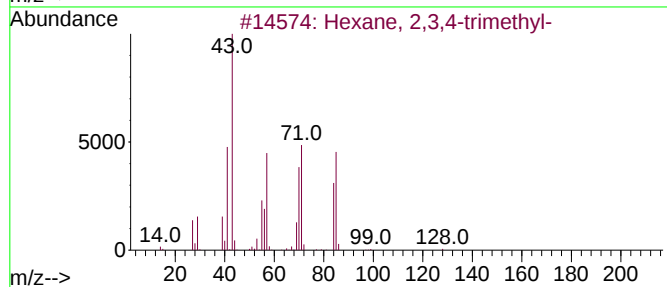
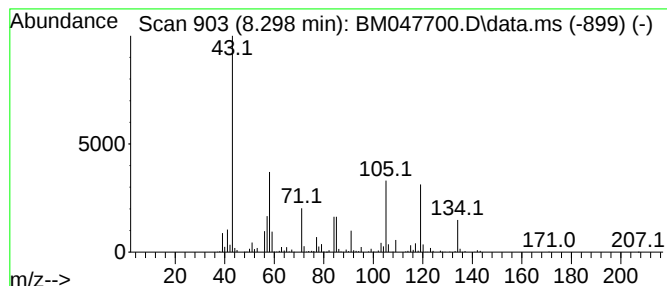
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.298	8.77 ng/ul	1544560	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	18
2			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	16
3			Octane, 4-methyl-	128	C9H20	002216-34-4	14
4			3-Oxetanol, 2,2,3-trimethyl-	116	C6H12O2	025910-96-7	12
5			2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

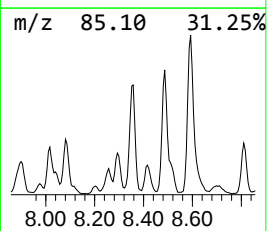
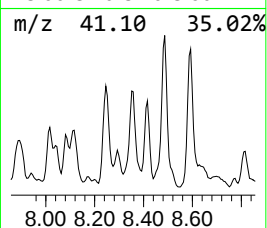
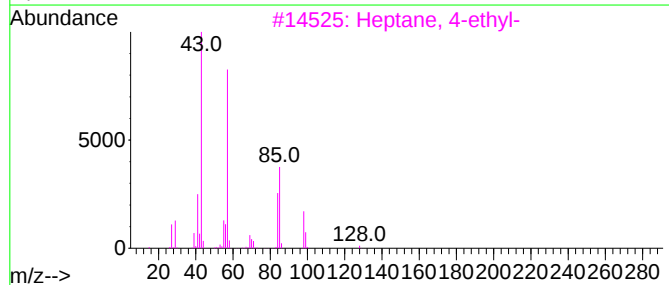
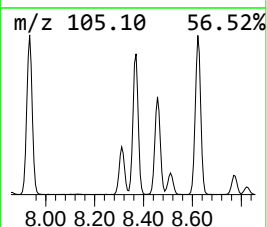
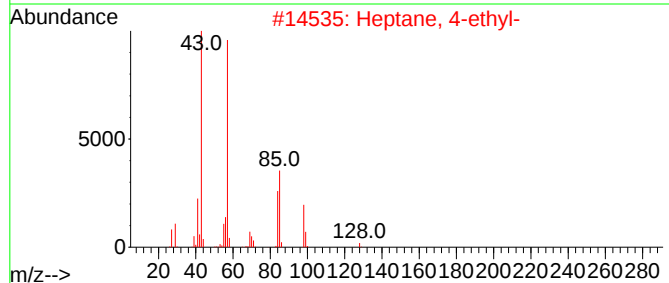
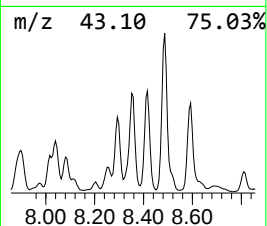
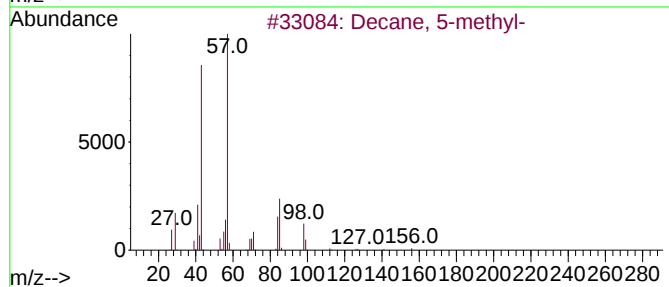
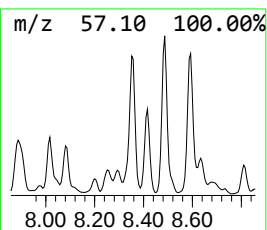
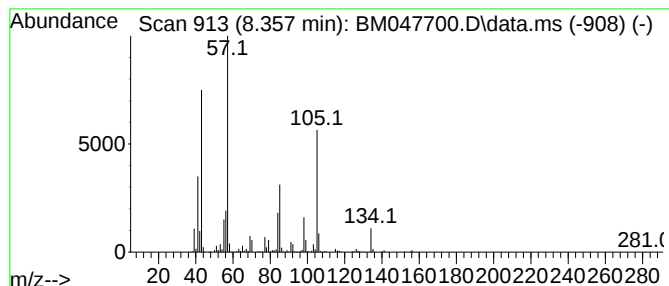
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.357	17.30 ng/ul	3046930	1,4-Dichlorobenzene-d4			7.816
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 5-methyl-		156	C11H24	013151-35-4	64
2	Heptane, 4-ethyl-		128	C9H20	002216-32-2	38
3	Heptane, 4-ethyl-		128	C9H20	002216-32-2	38
4	2,2-Dimethylpropionic acid, 4-me...		186	C11H22O2	150588-62-8	27
5	Octane, 2,3-dimethyl-		142	C10H22	007146-60-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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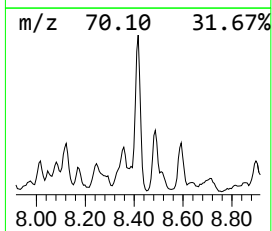
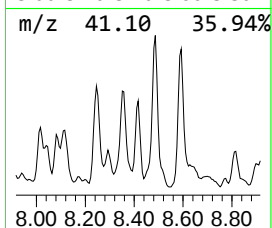
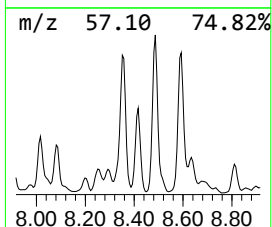
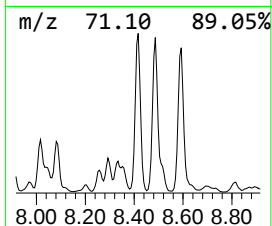
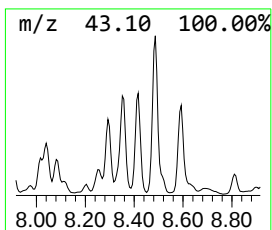
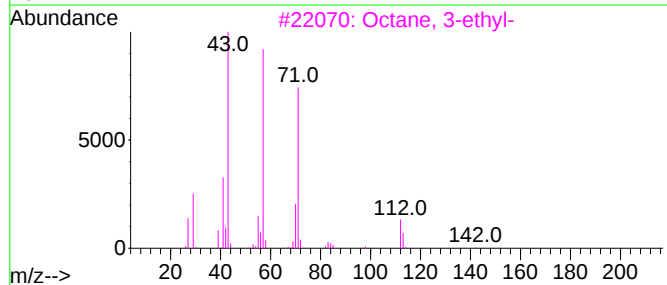
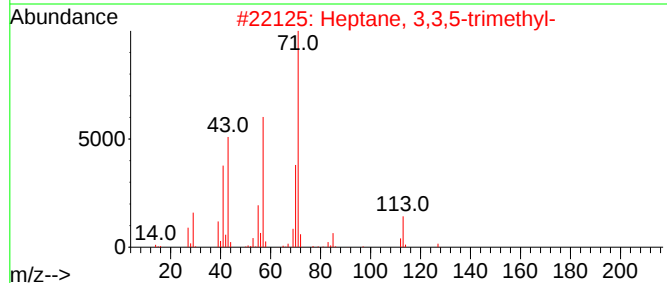
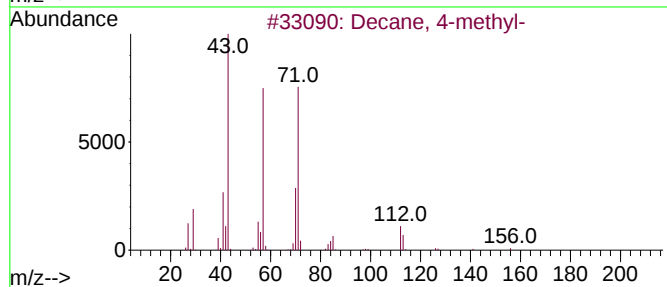
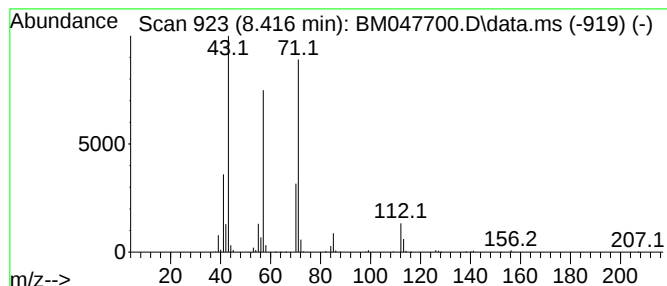
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.416	8.61 ng/ul	1516470	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	91
2			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
3			Octane, 3-ethyl-	142	C10H22	005881-17-4	78
4			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	72
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
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Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

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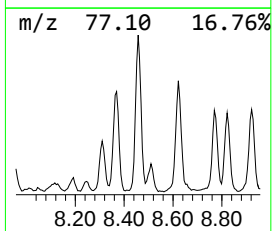
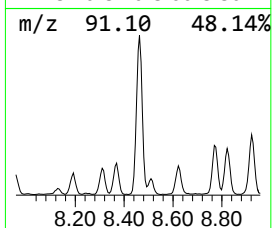
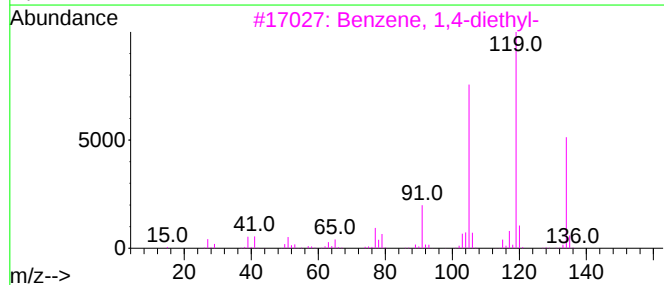
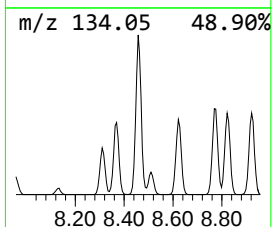
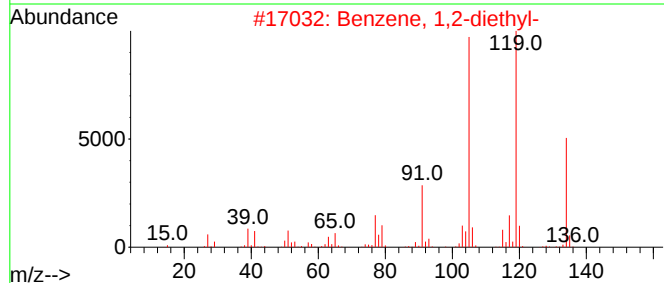
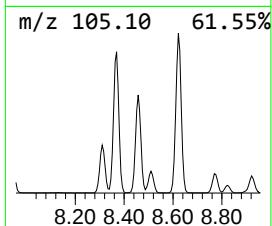
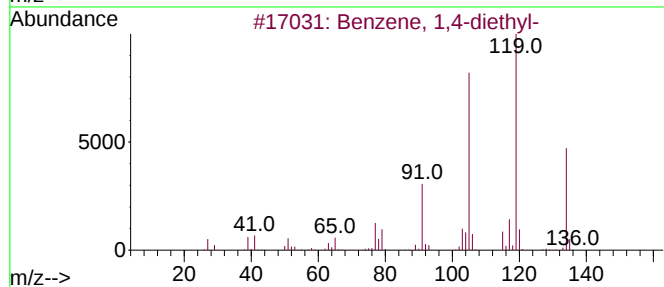
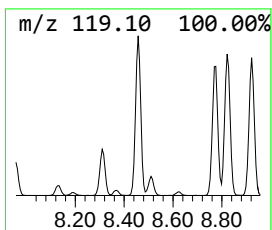
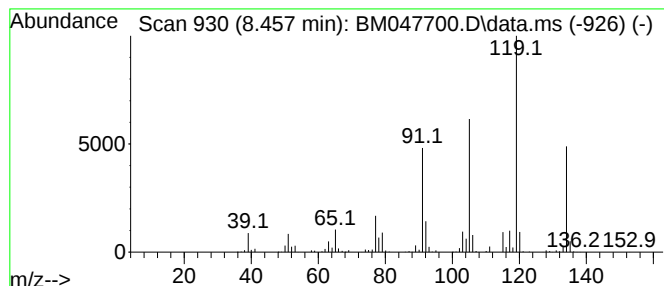
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzene, 1,4-diethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.457	11.43 ng/ul	2013230	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
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Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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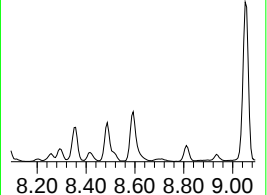
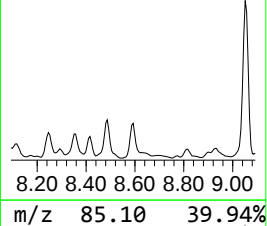
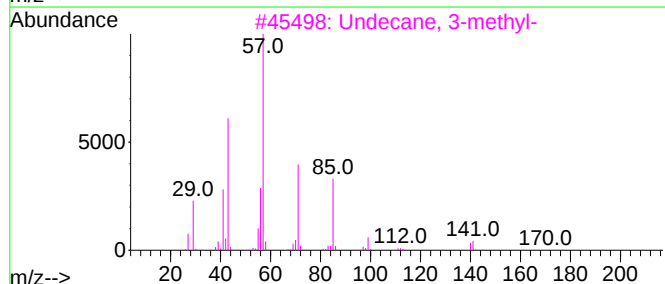
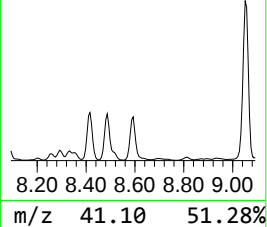
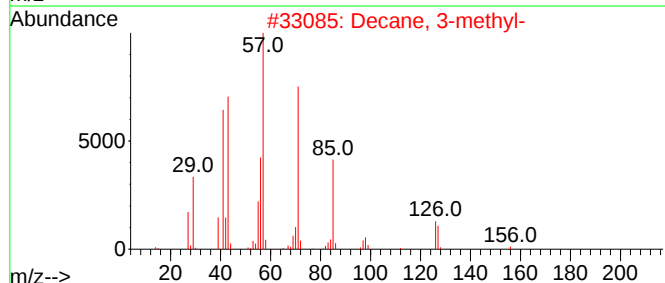
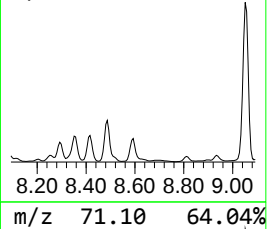
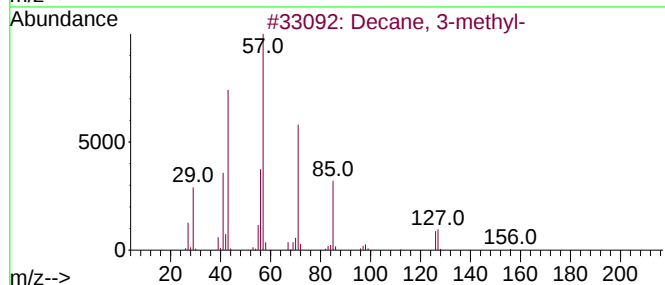
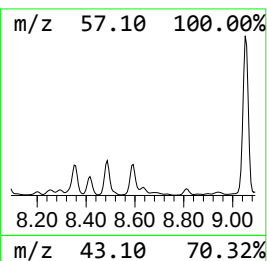
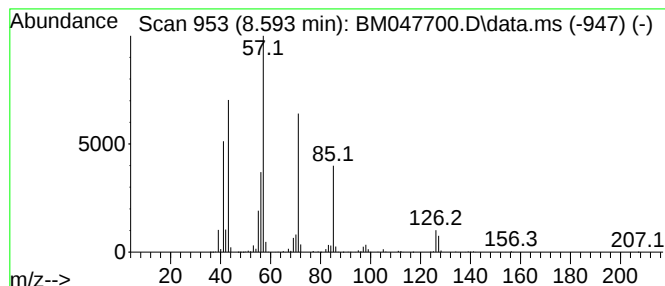
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.593	14.44 ng/ul	2541780	1,4-Dichlorobenzene-d4	7.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 3-methyl-	156	C11H24	013151-34-3	81
2		Decane, 3-methyl-	156	C11H24	013151-34-3	78
3		Undecane, 3-methyl-	170	C12H26	001002-43-3	64
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
5		1-Iodoundecane	282	C11H23I	004282-44-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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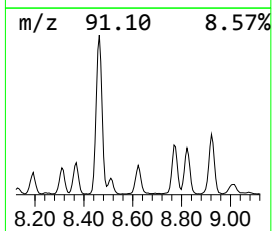
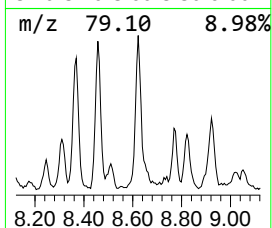
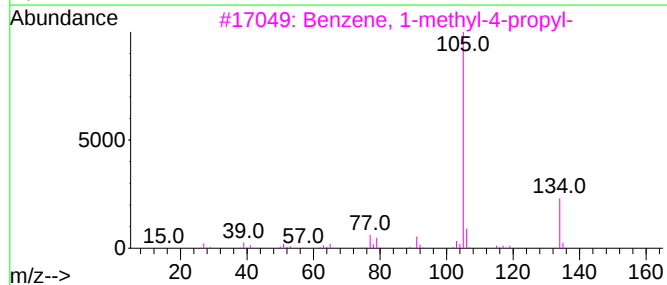
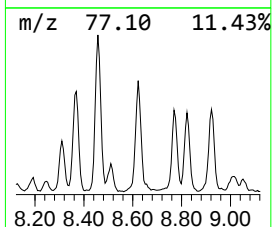
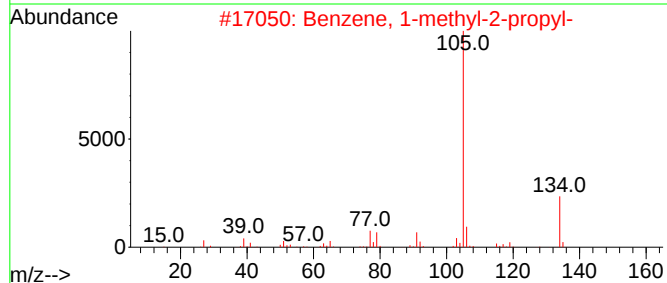
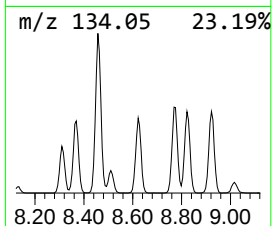
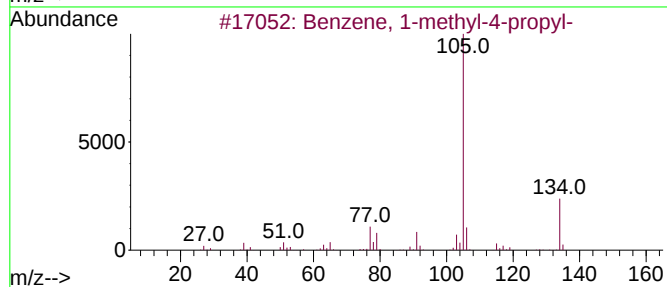
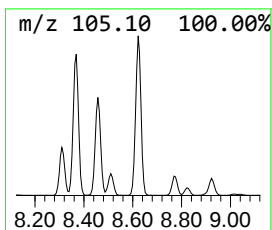
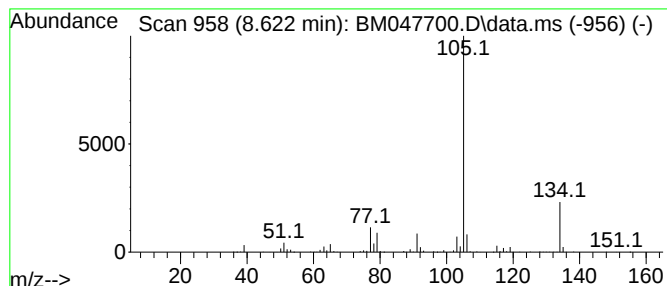
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 1-methyl-4-propyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.622	7.16 ng/ul	1261080	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

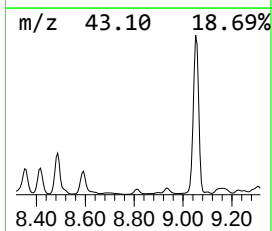
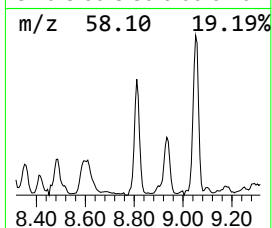
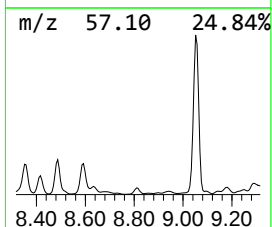
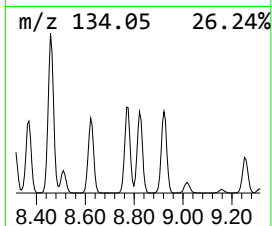
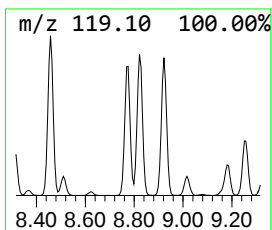
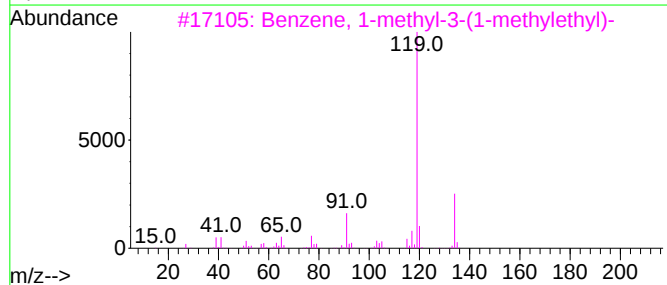
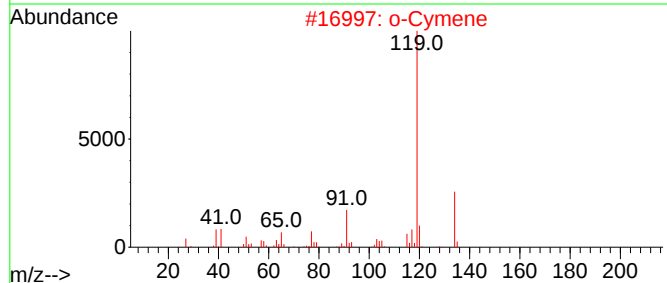
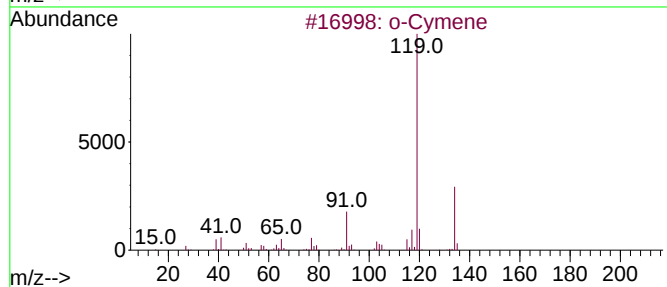
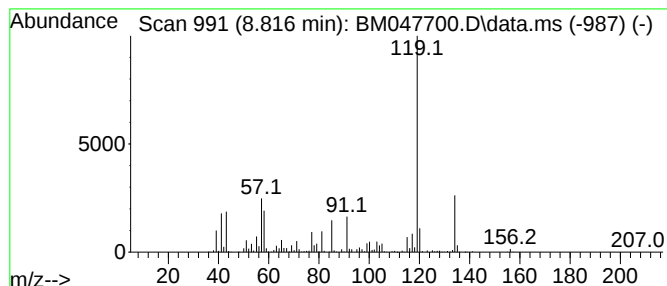
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 o-Cymene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.816	10.08 ng/ul	1774830	1,4-Dichlorobenzene-d4			7.816
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	o-Cymene		134	C10H14	000527-84-4	95
2	o-Cymene		134	C10H14	000527-84-4	95
3	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	95
4	p-Cymene		134	C10H14	000099-87-6	93
5	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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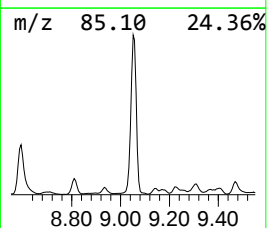
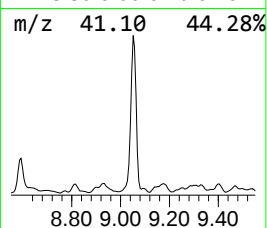
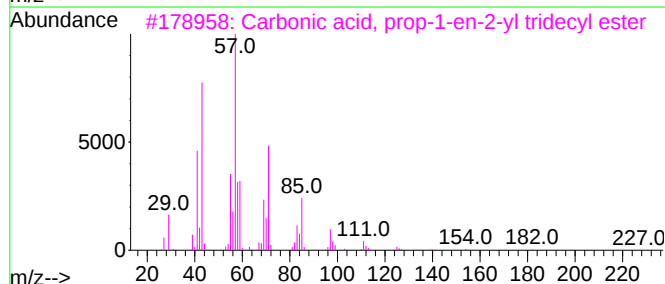
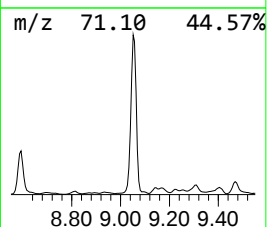
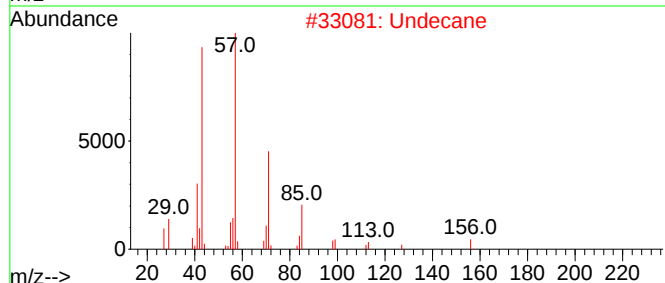
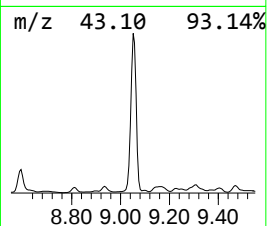
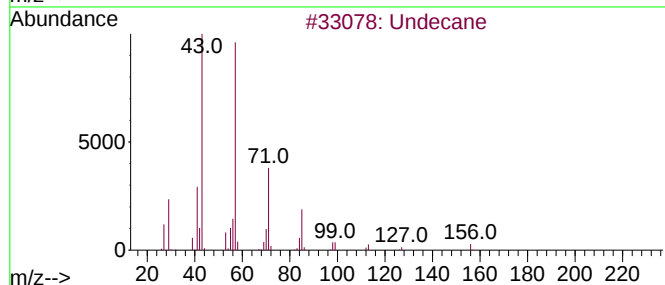
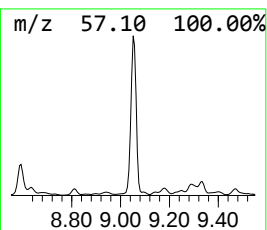
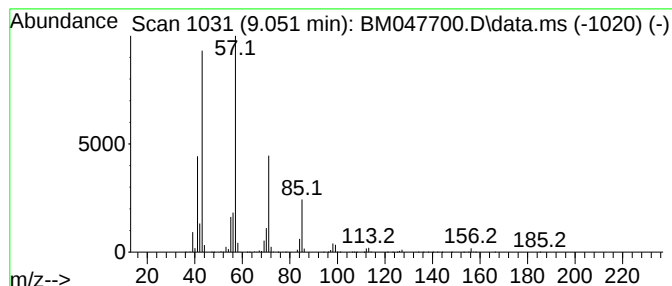
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.051	63.89 ng/ul	11249500	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Undecane	156	C11H24	001120-21-4	91
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	90
4			Tridecane	184	C13H28	000629-50-5	86
5			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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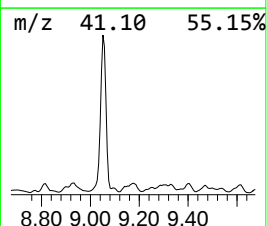
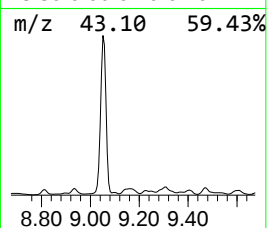
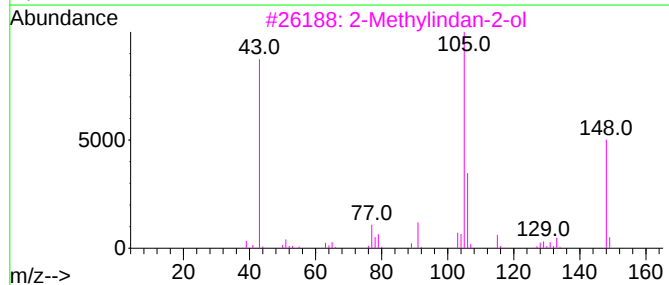
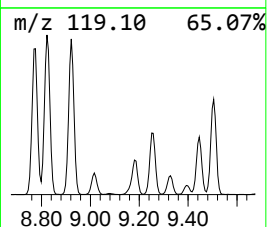
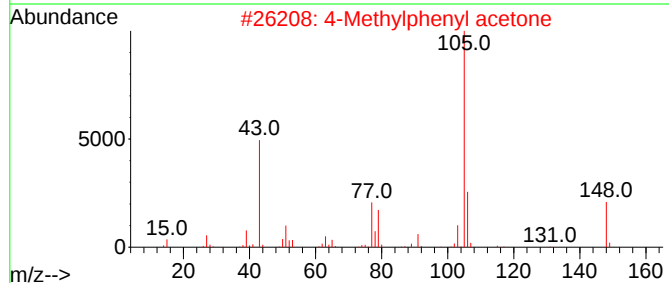
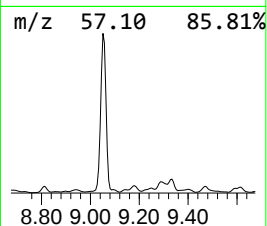
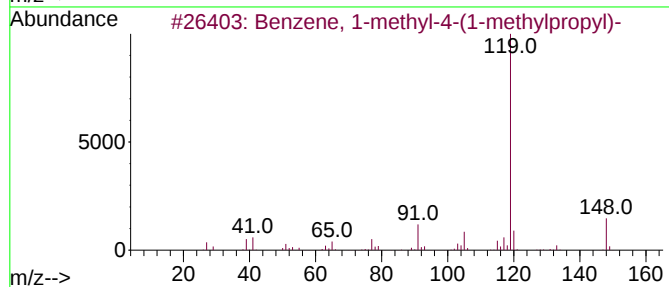
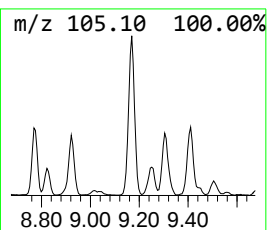
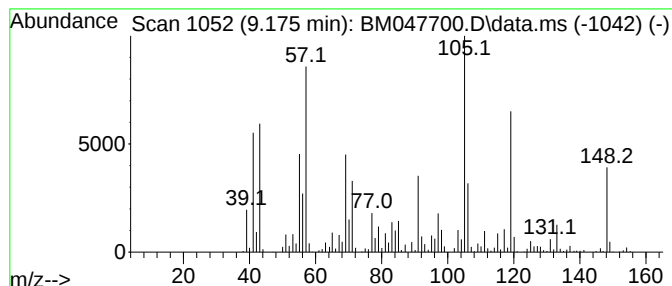
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Benzene, 1-methyl-4-(1-meth... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.175	12.11 ng/ul	2133140	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	55
2			4-Methylphenyl acetone	148	C10H12O	002096-86-8	50
3			2-Methylindan-2-ol	148	C10H12O	033223-84-6	45
4			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	43
5			2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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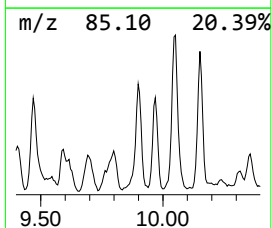
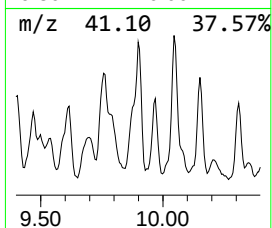
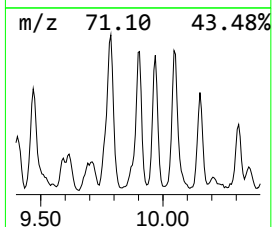
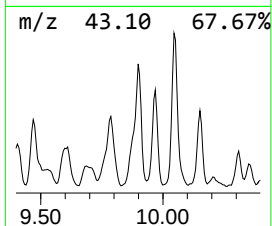
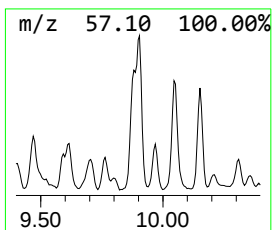
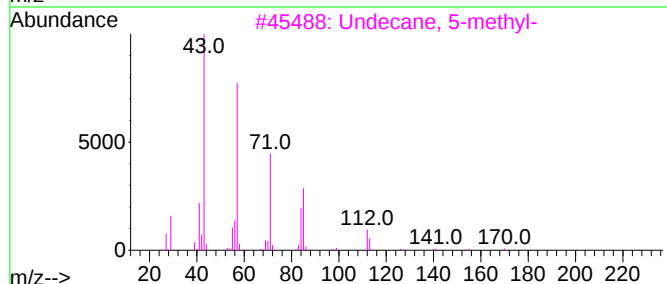
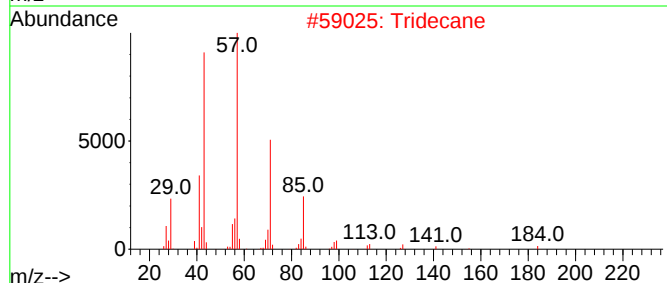
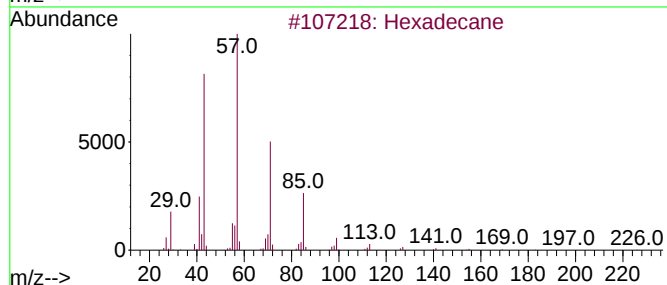
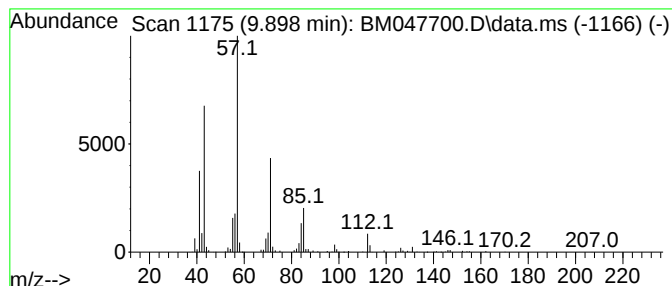
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.898	2.43 ng/ul	1907730	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	80
2			Tridecane	184	C13H28	000629-50-5	72
3			Undecane, 5-methyl-	170	C12H26	001632-70-8	59
4			Nonane, 5-(1-methylpropyl)-	184	C13H28	062185-54-0	59
5			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
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Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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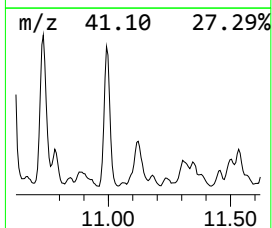
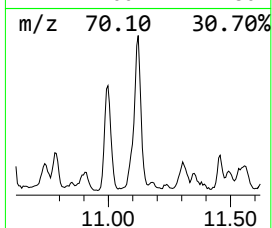
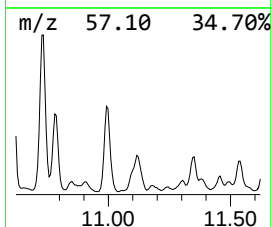
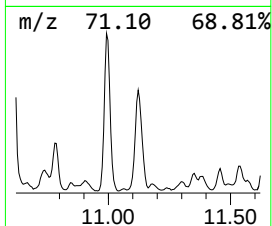
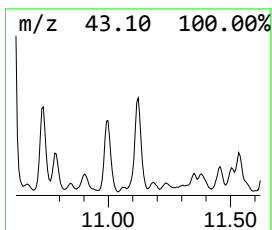
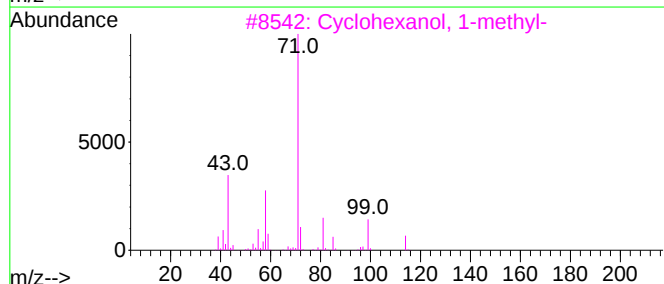
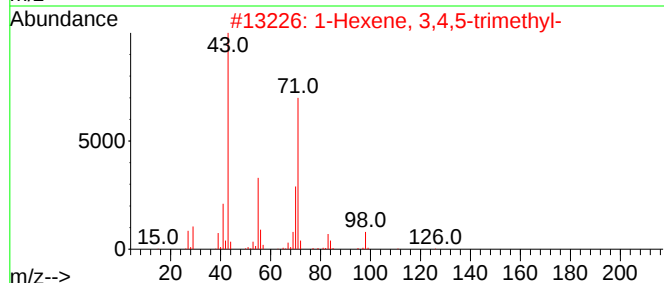
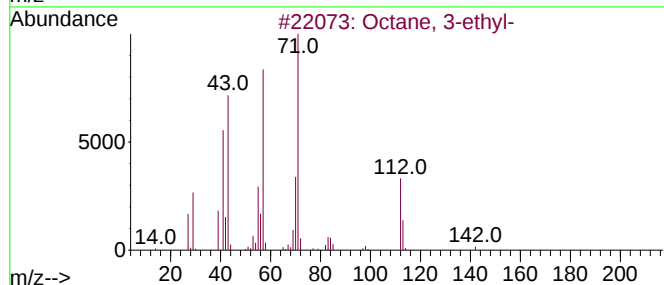
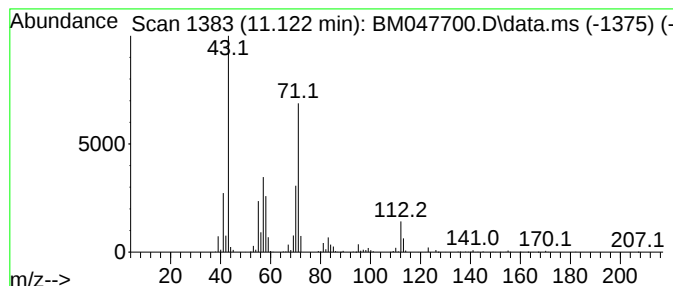
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.122	2.13 ng/ul	1672730	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-			142	C10H22	005881-17-4	50
2	1-Hexene, 3,4,5-trimethyl-			126	C9H18	056728-10-0	47
3	Cyclohexanol, 1-methyl-			114	C7H14O	000590-67-0	46
4	Carbonic acid, neopentyl 2-ethyl...			244	C14H28O3	1000357-85-7	43
5	Decane, 4-methyl-			156	C11H24	002847-72-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
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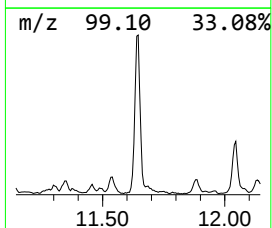
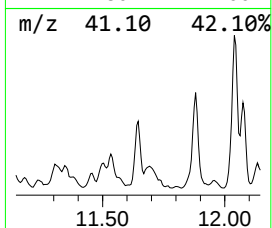
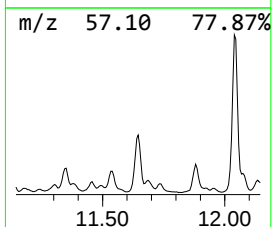
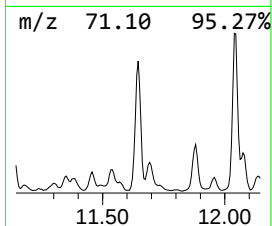
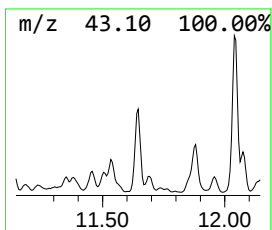
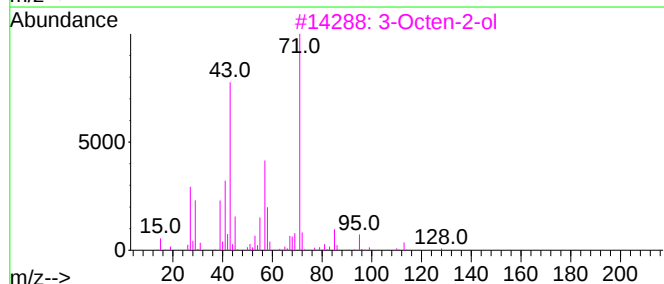
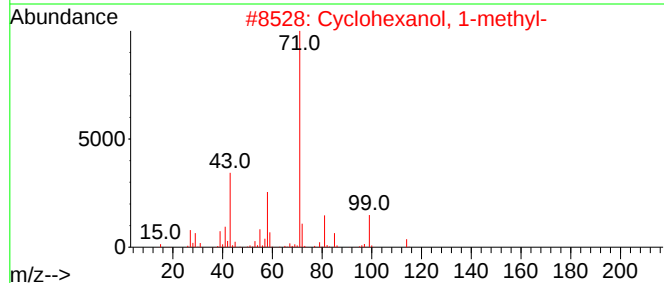
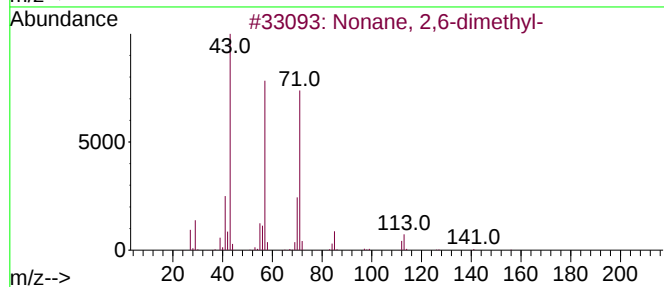
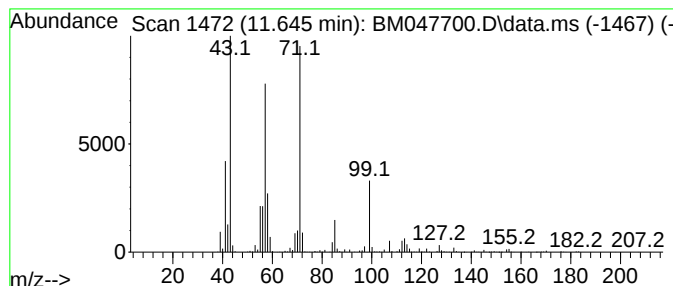
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.645	2.54 ng/ul	2000270	Naphthalene-d8	10.610	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	58
2	Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	52
3	3-Octen-2-ol	128	C8H16O	076649-14-4	52
4	Octane, 2-bromo-	192	C8H17Br	000557-35-7	47
5	Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

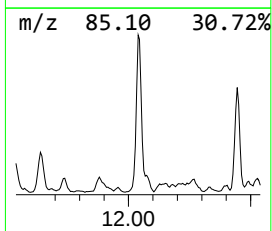
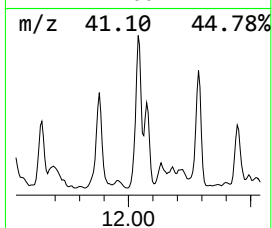
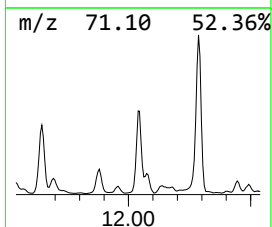
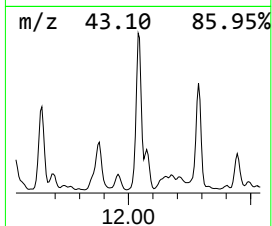
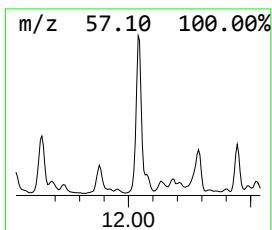
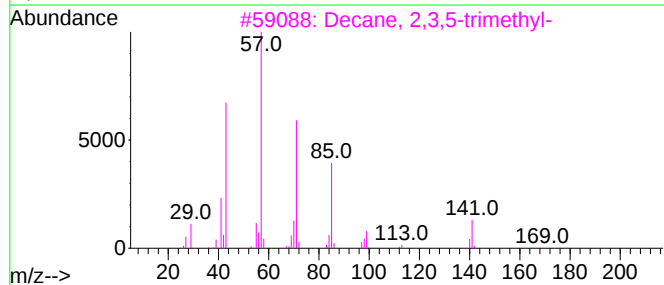
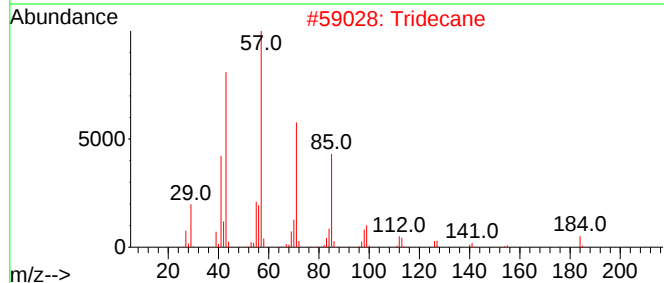
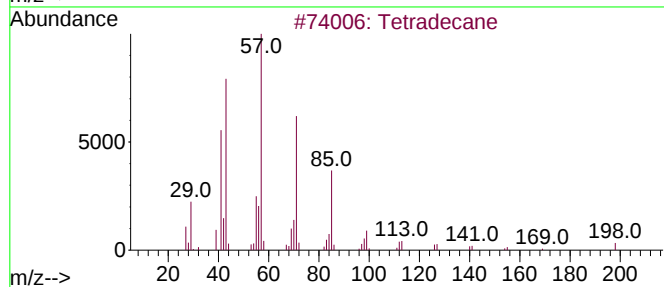
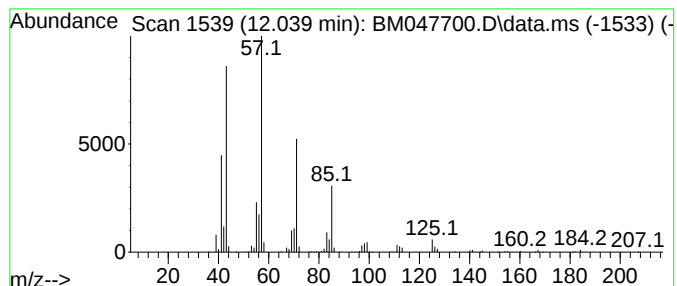
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ClientSampleId :
DDC93DL

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Quant Title : SVOA CALIBRATION

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

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.039	4.55 ng/ul	3577930	Naphthalene-d8	10.610	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	80
2	Tridecane	184	C13H28	000629-50-5	80
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	80
4	Dodecane	170	C12H26	000112-40-3	80
5	Tridecane	184	C13H28	000629-50-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

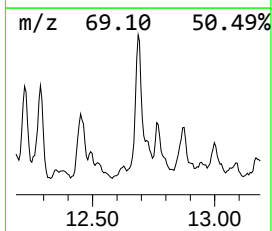
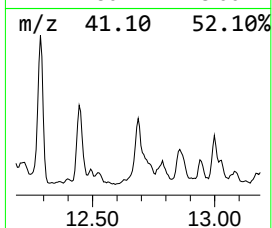
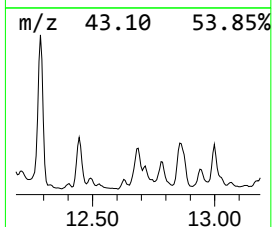
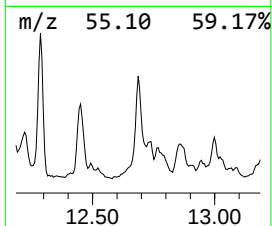
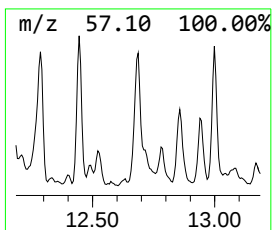
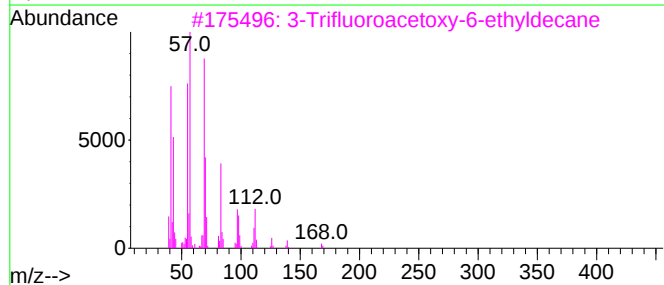
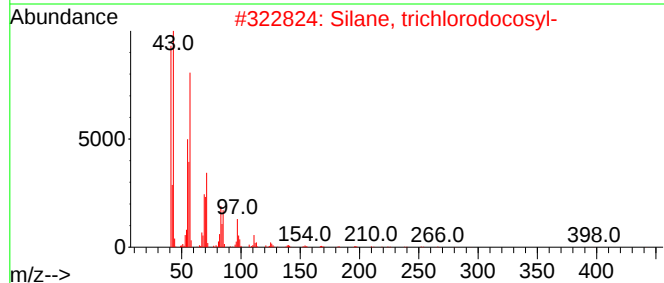
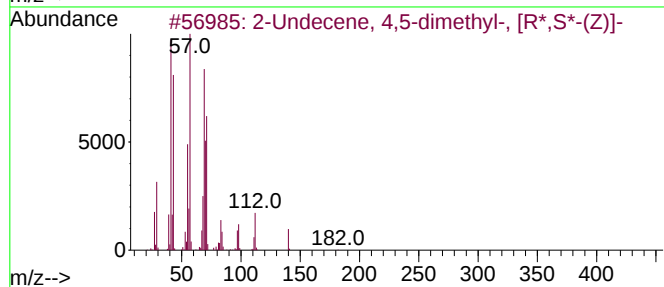
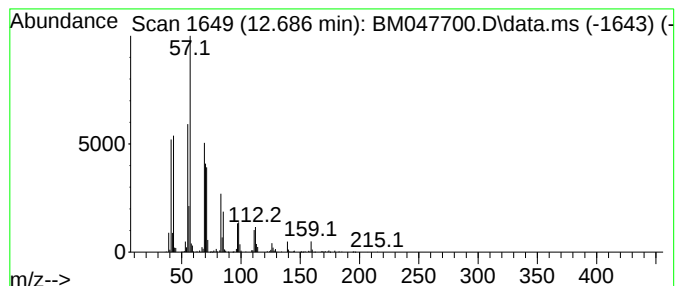
Instrument :
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ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.686	14.25 ng/ul	1800050	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Undecene, 4,5-dimethyl-, [R*,S...		182	C13H26	055170-93-9	58
2	Silane, trichlorodocosyl-		442	C22H45Cl3Si	007325-84-0	53
3	3-Trifluoroacetoxy-6-ethyldecane		282	C14H25F3O2	116436-59-0	46
4	Heptane, 1,1'-oxybis-		214	C14H30O	000629-64-1	43
5	1-Hexene, 3,3-dimethyl-		112	C8H16	003404-77-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DDC93DL

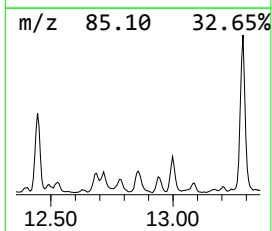
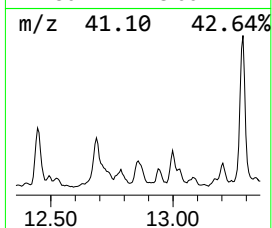
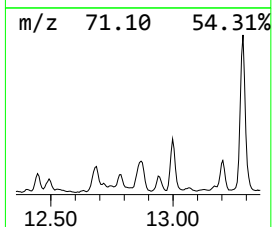
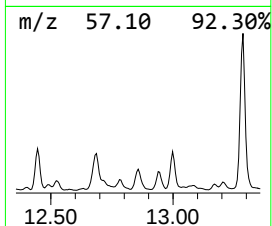
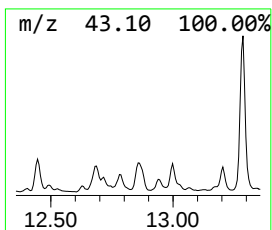
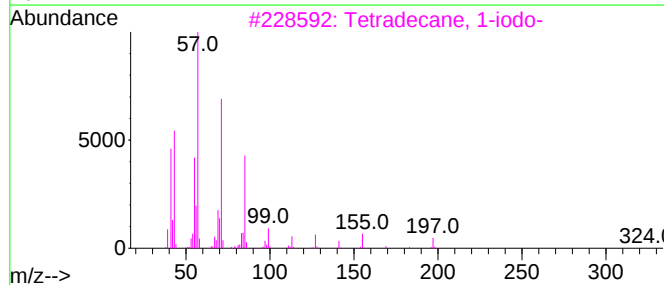
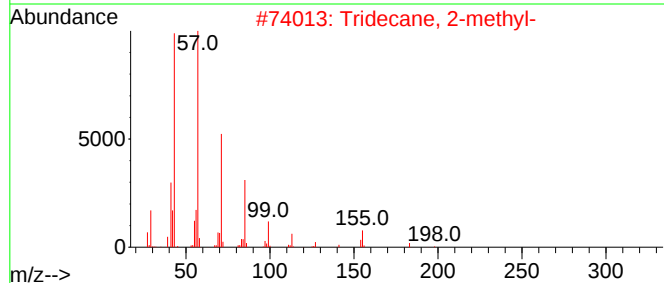
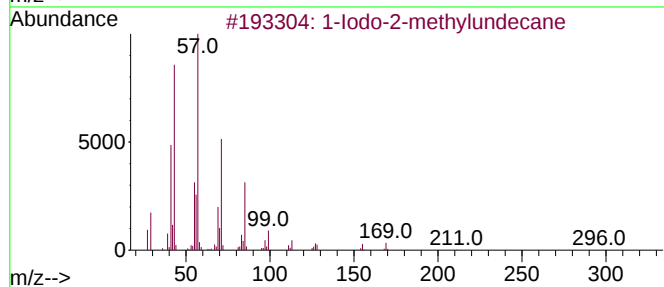
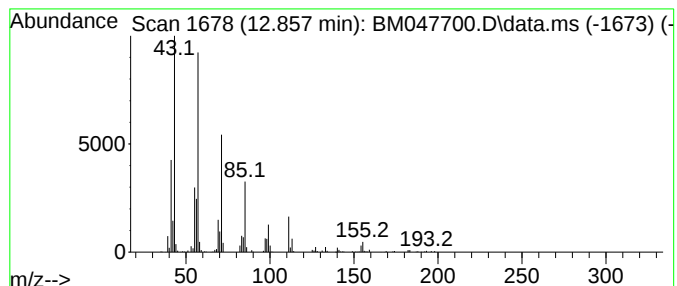
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 1-Iodo-2-methylundecane Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.857	6.80 ng/ul	858165	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	86
2			Tridecane, 2-methyl-	198	C14H30	001560-96-9	76
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

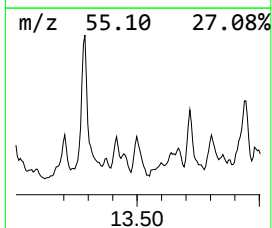
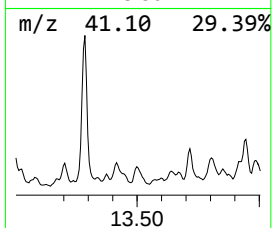
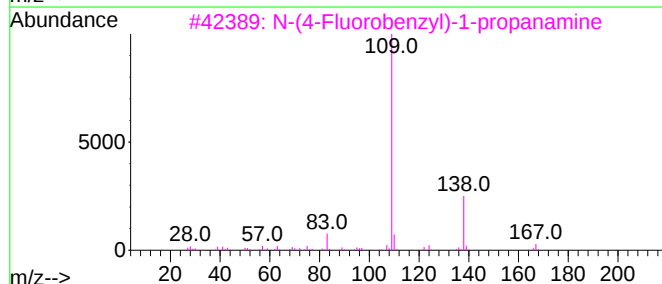
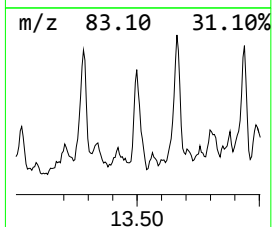
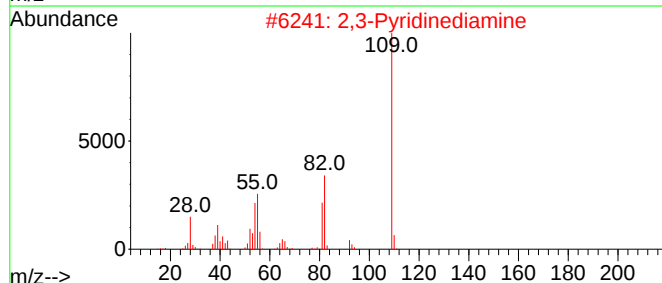
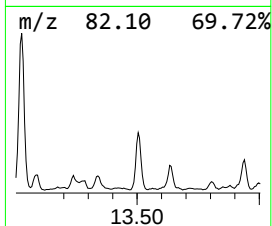
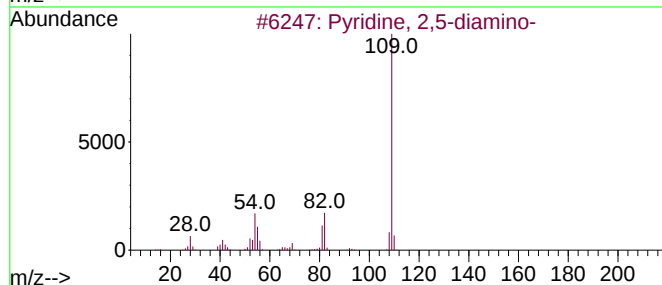
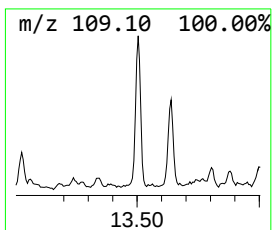
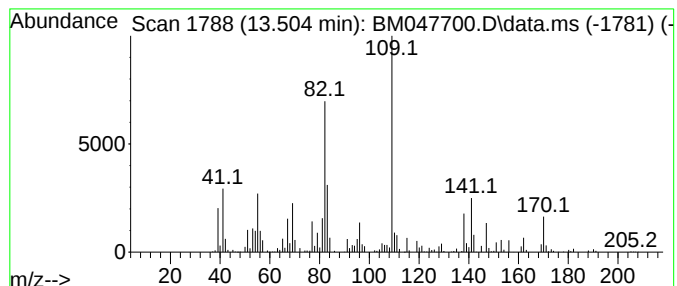
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 unknown-02

Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.504	10.79 ng/ul	1362300	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 2,5-diamino-	109	C5H7N3	004318-76-7	47
2			2,3-Pyridinediamine	109	C5H7N3	000452-58-4	47
3			N-(4-Fluorobenzyl)-1-propanamine	167	C10H14FN	741698-80-6	43
4			2-(5-Methyl-furan-2-yl)-propiona...	138	C8H10O2	1000193-72-3	41
5			2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
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ClientSampleId :
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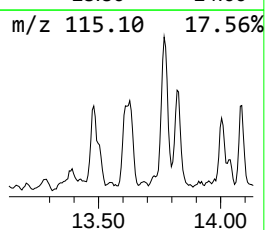
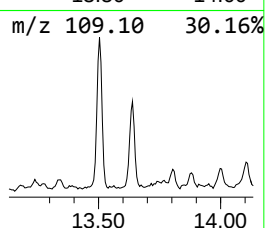
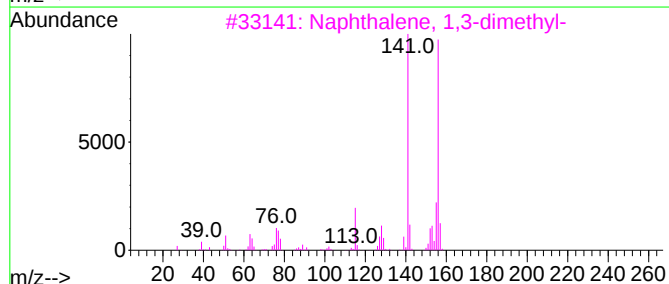
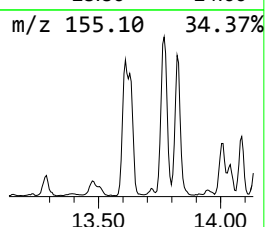
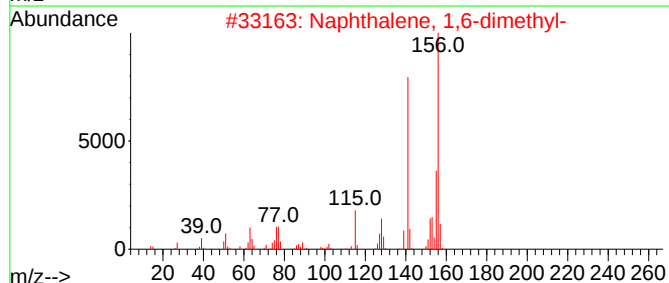
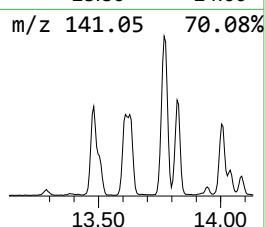
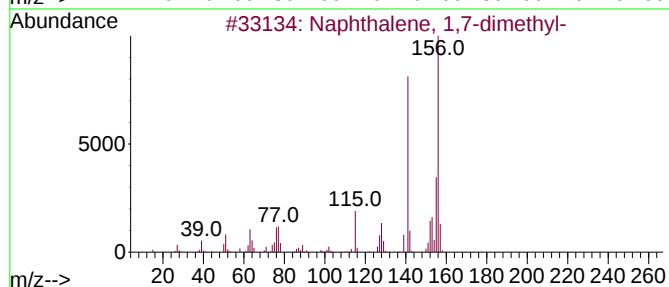
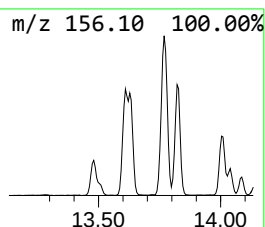
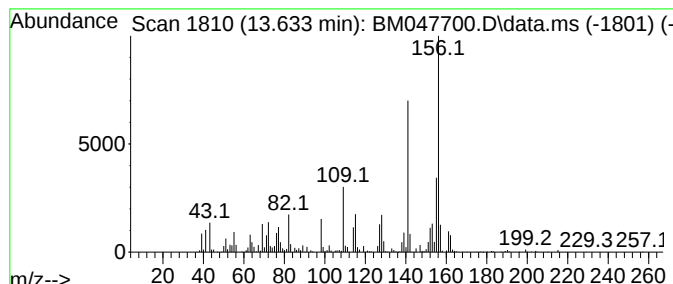
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Naphthalene, 1,7-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.633	15.78 ng/ul	1992370	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

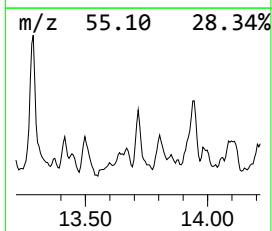
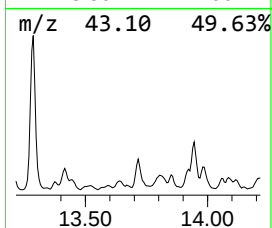
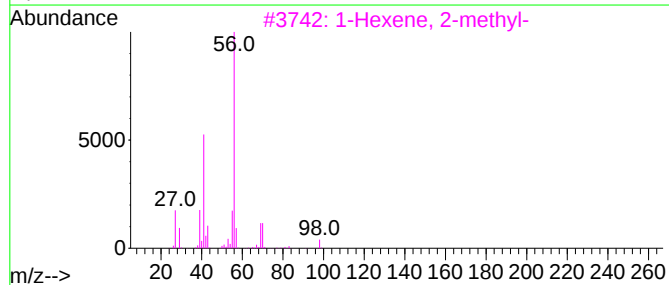
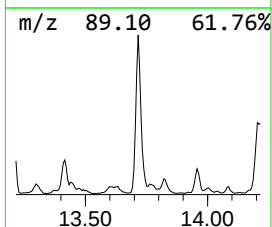
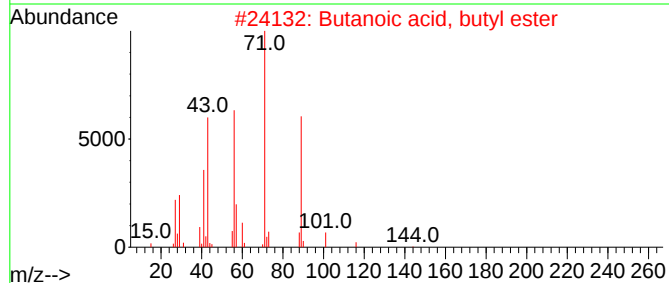
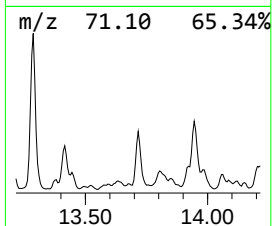
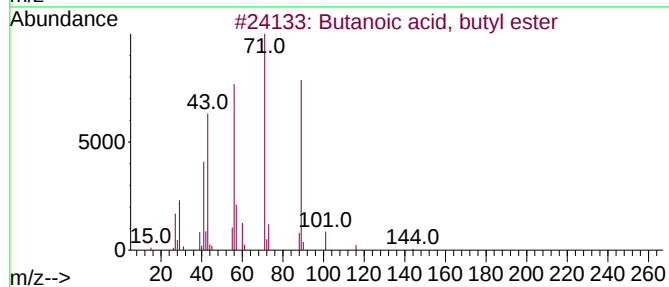
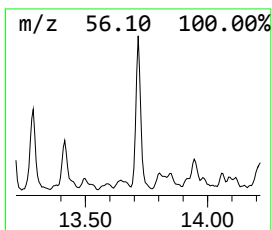
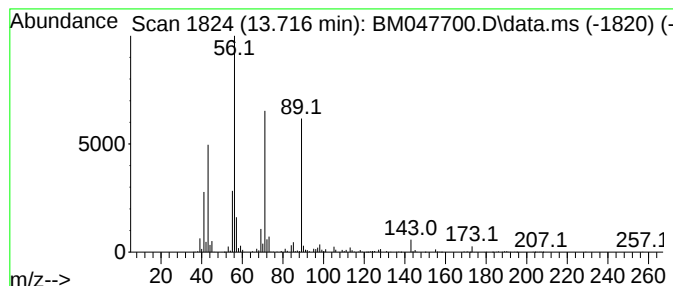
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 Butanoic acid, butyl ester Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	8.04 ng/ul	1015040	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	52
2			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	50
3			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
4			Nonane, 5-methylene-	140	C10H20	006795-79-5	38
5			3,4,5-Trimethyldihydrofuran-2-one	128	C7H12O2	1000194-05-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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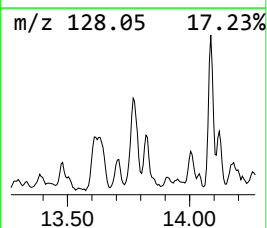
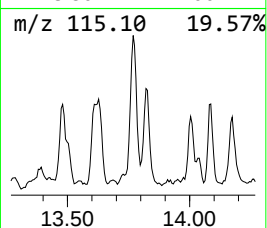
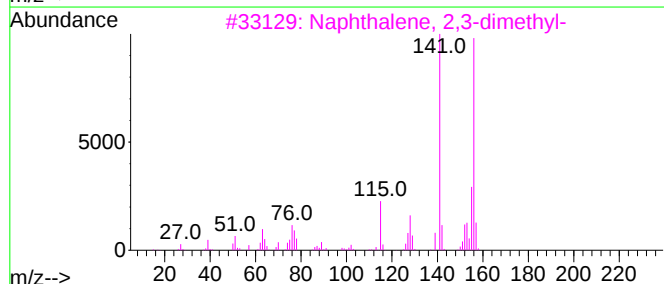
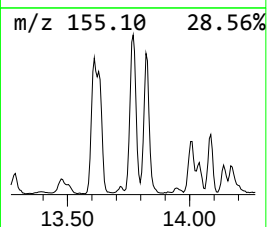
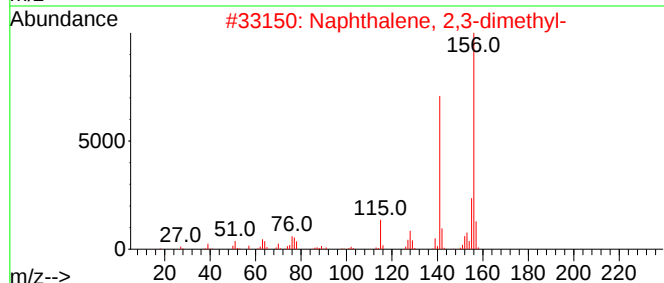
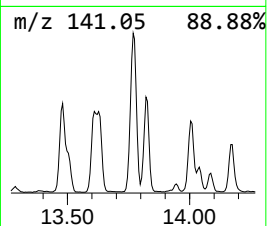
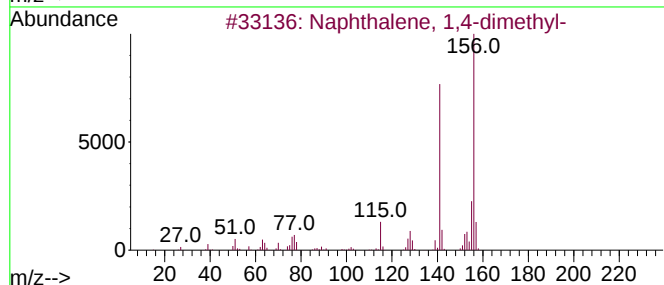
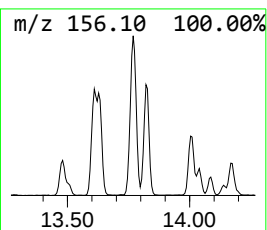
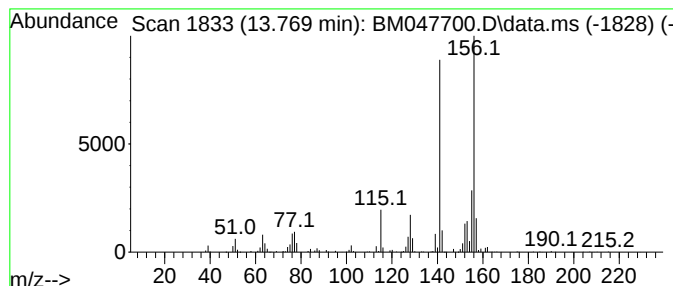
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Naphthalene, 1,4-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	12.27 ng/ul	1549120	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

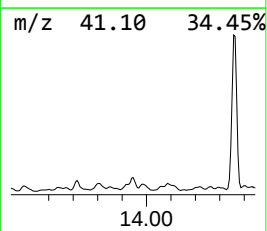
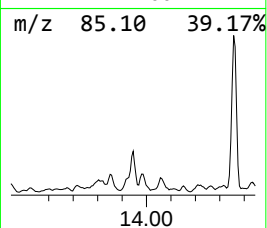
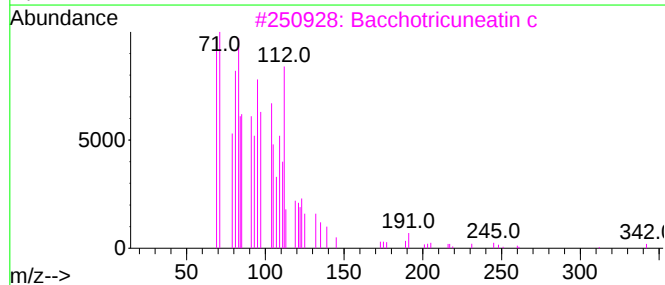
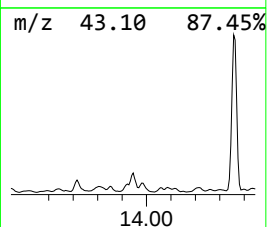
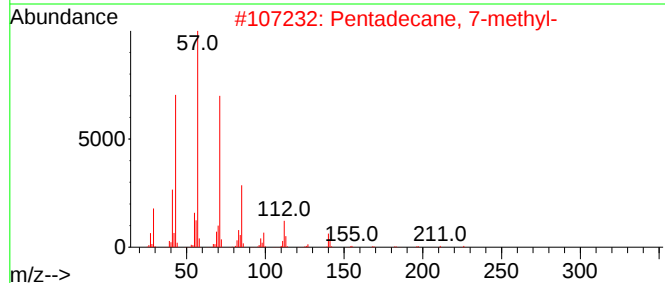
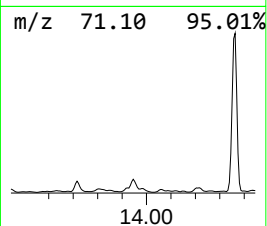
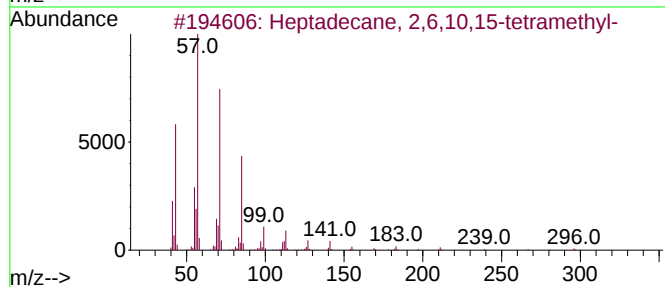
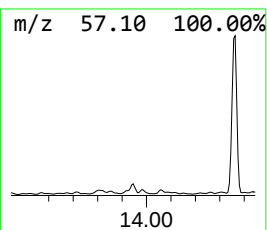
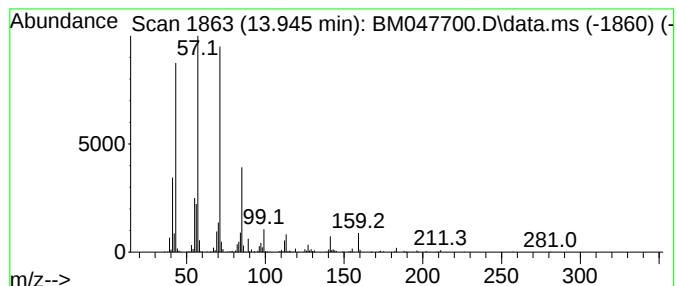
Instrument :
BNA_M
ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.945	8.24 ng/ul	1040840	Acenaphthene-d10			14.445
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	86
2	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	78
3	Bacchotricuneatin c		342	C20H22O5	066563-30-2	72
4	Hexadecane		226	C16H34	000544-76-3	64
5	Tridecane, 5-propyl-		226	C16H34	055045-11-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

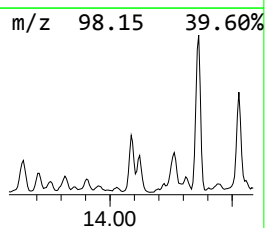
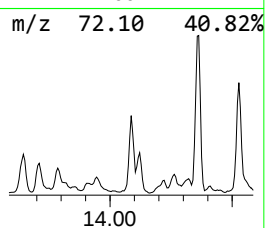
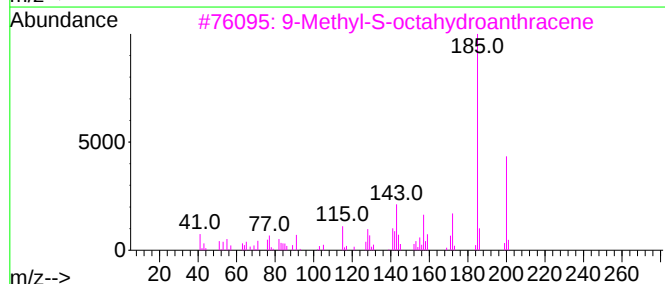
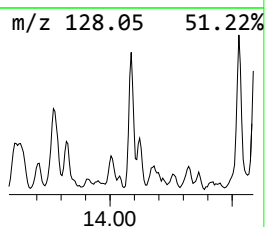
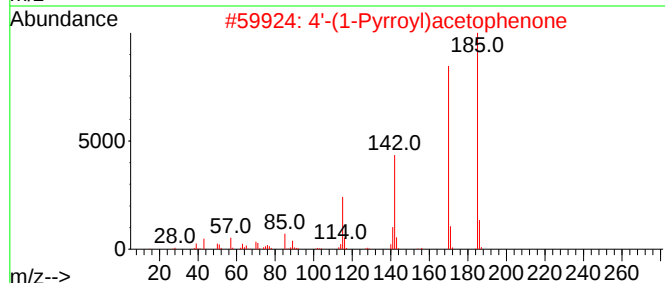
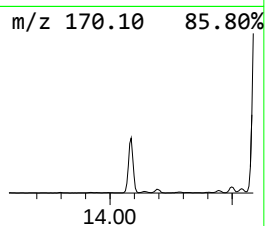
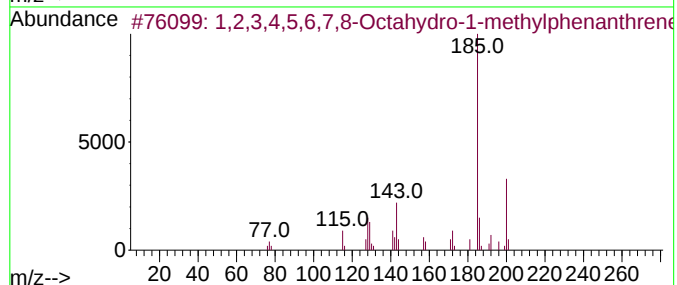
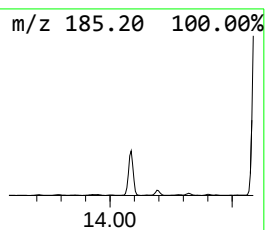
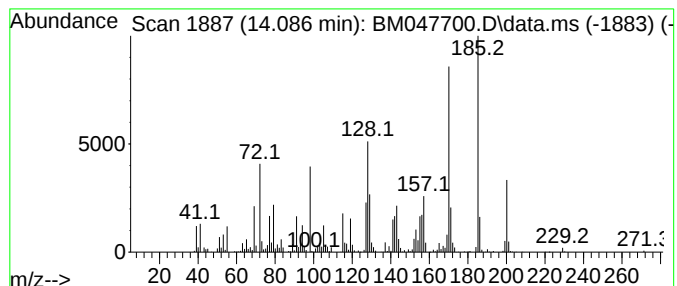
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 unknown-03 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	13.95 ng/ul	1761470	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	42		
2	4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	38		
3	9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	38		
4	Silane, trimethyl-2-naphthalenyl-	200	C13H16Si	018052-85-2	35		
5	Ethanone, 1-(5-methyl-1-phenyl-1...	200	C12H12N2O	006123-63-3	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

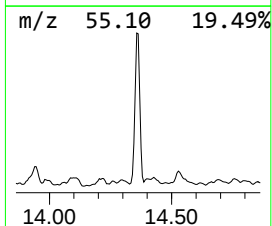
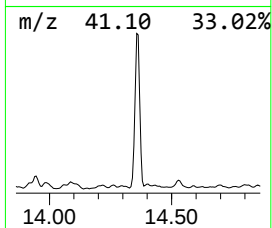
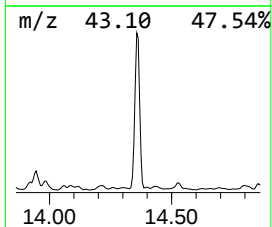
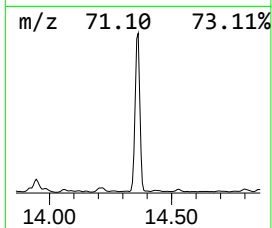
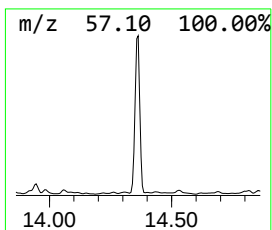
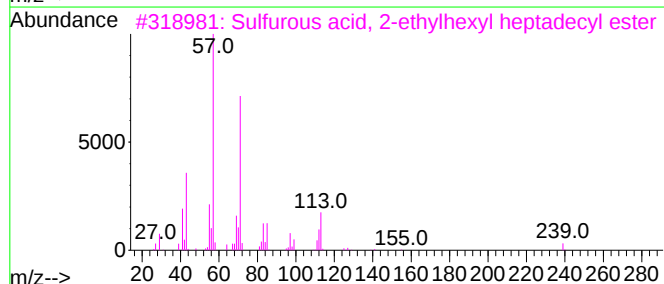
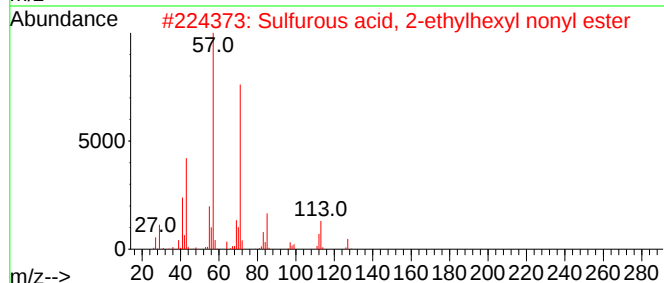
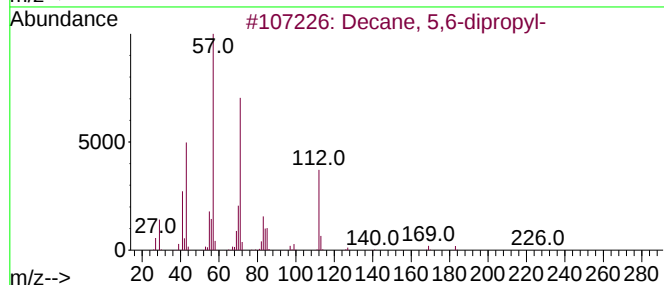
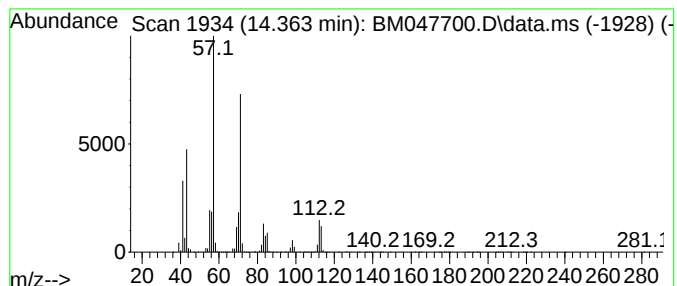
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.363	104.30 ng/ul	13171600	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	72
2	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72
3	Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	72
4	2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	72
5	Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DDC93DL

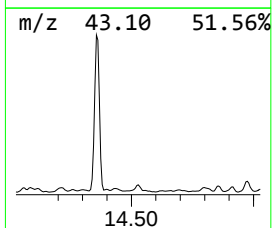
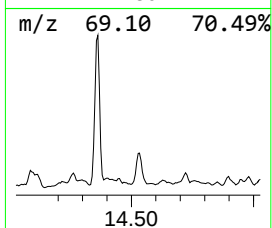
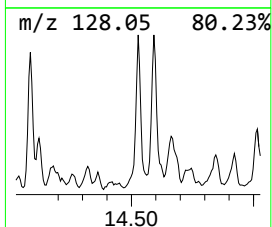
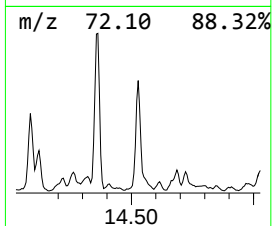
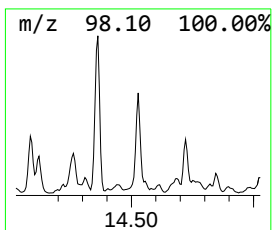
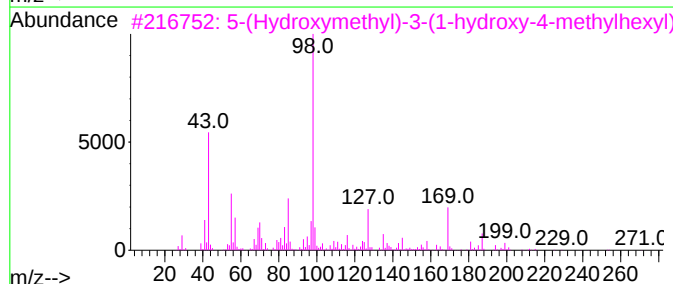
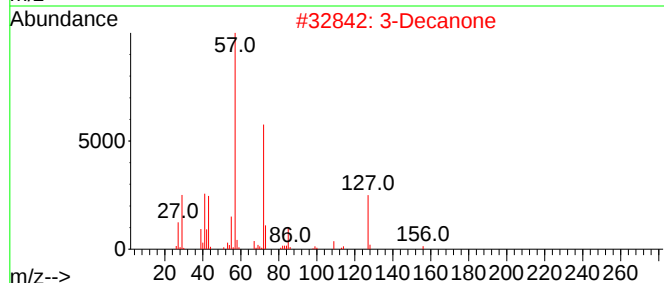
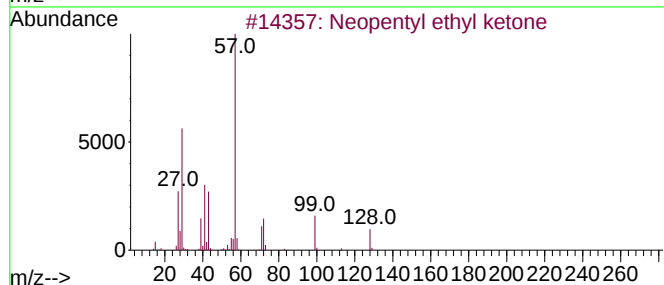
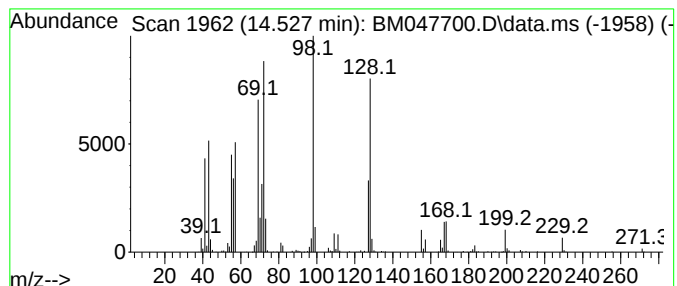
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 unknown-04 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.527	8.72 ng/ul	1101760	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Neopentyl ethyl ketone	128	C8H16O	005340-30-7	38
2			3-Decanone	156	C10H20O	000928-80-3	22
3			5-(Hydroxymethyl)-3-(1-hydroxy-4...	314	C16H26O6	1000480-56-8	22
4			Carbaminic acid,2-(4-chloropheno...	258	C10H11ClN2O4	1010140-52-2	14
5			2-(2-Chlorophenoxy)ethanol, isop...	214	C11H15ClO2	1000394-99-0	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

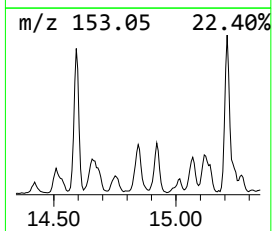
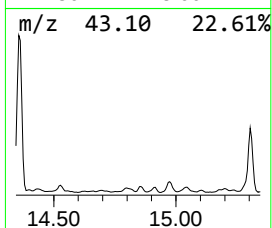
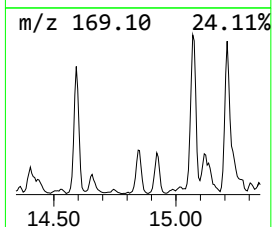
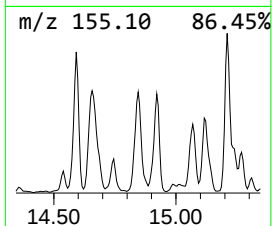
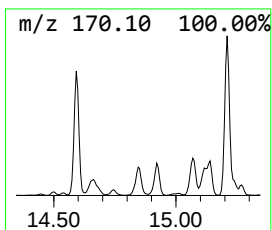
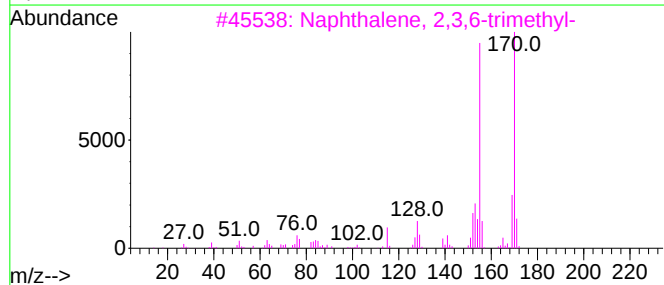
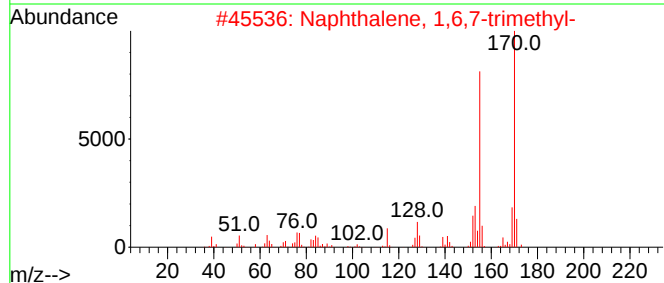
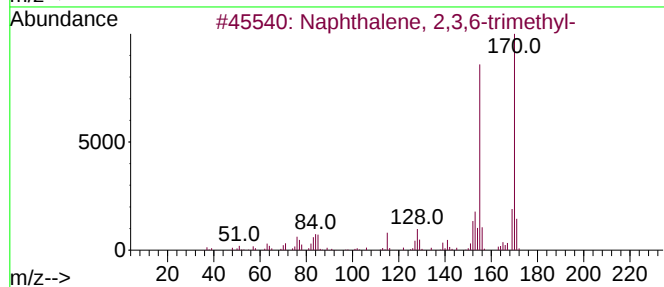
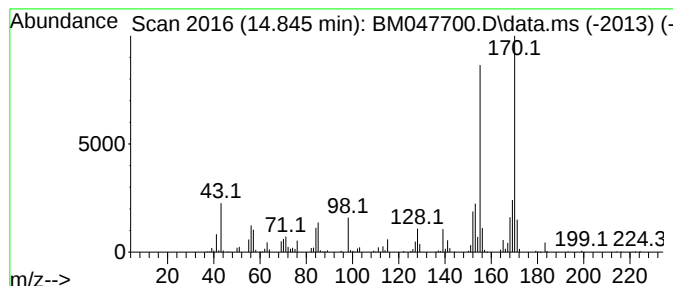
Instrument :
BNA_M
ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 Naphthalene, 2,3,6-trimethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.845	7.78 ng/ul	982753	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
4	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	91
5	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

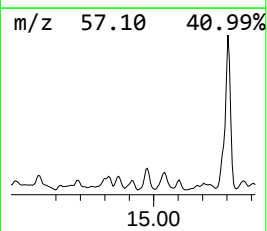
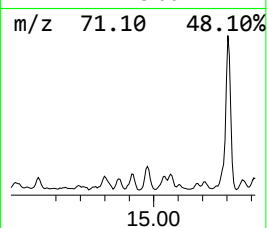
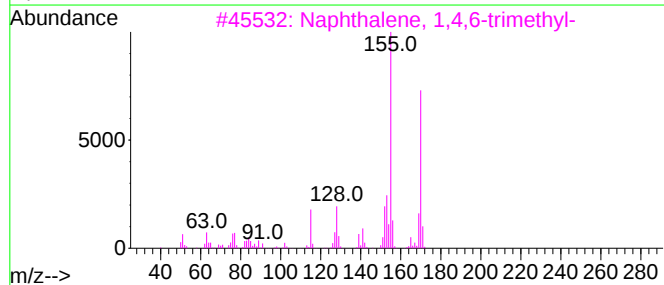
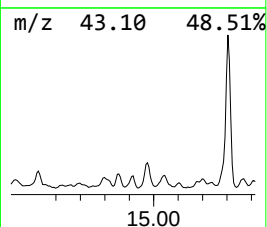
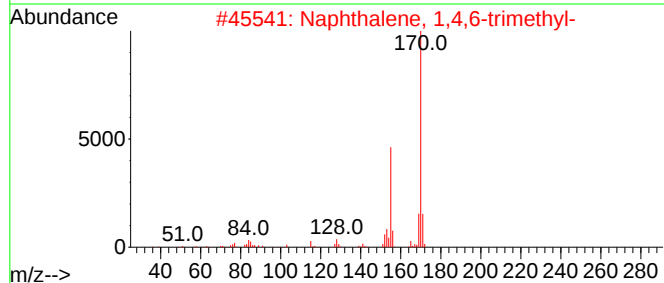
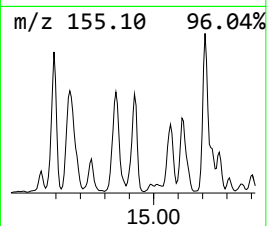
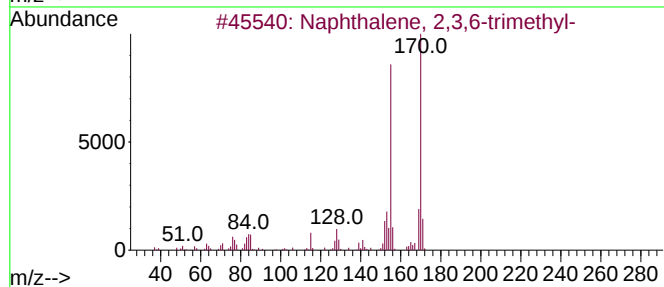
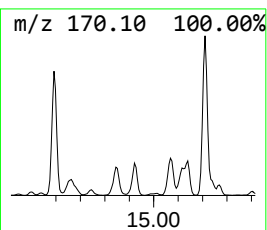
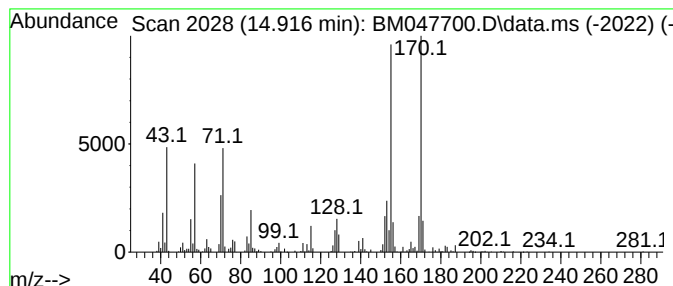
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 47 Naphthalene, 1,4,6-trimethyl- Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.916	6.95 ng/ul	878188	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

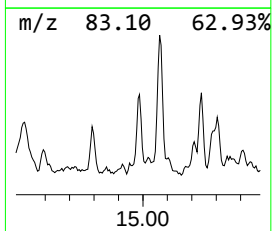
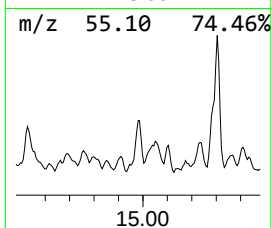
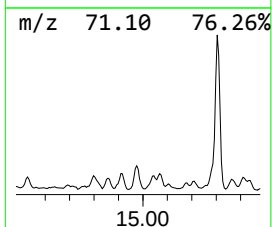
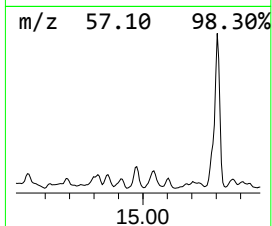
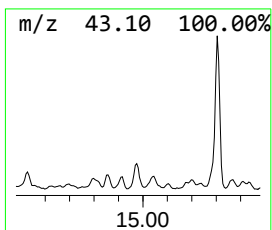
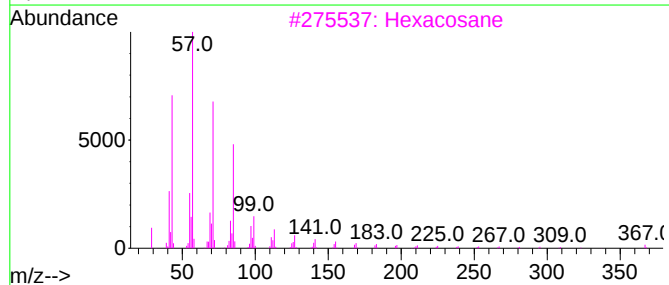
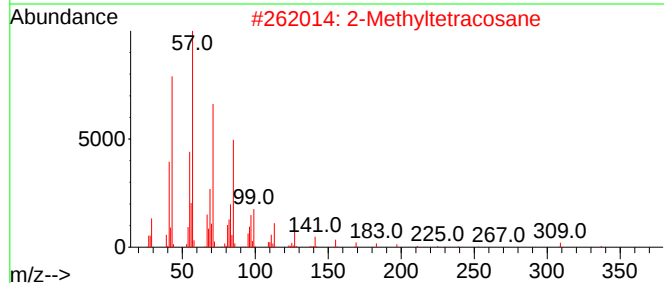
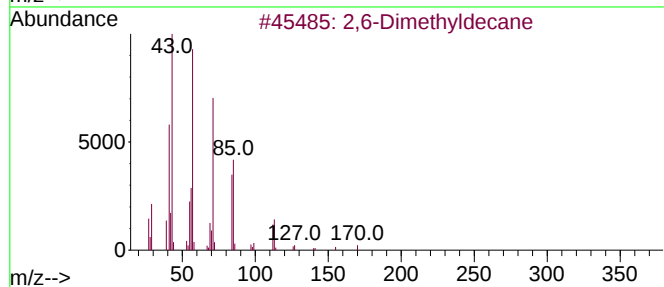
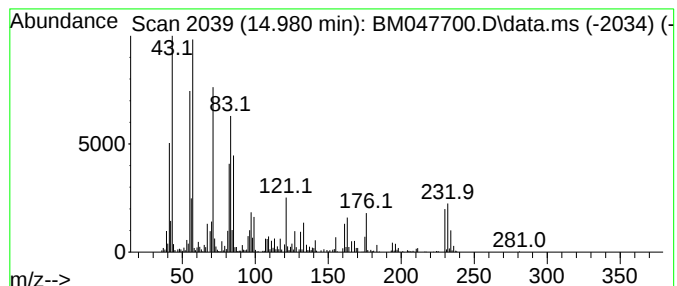
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ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.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 47

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.980	7.47 ng/ul	942715	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,6-Dimethyldecane		170	C12H26	013150-81-7	46
2	2-Methyltetracosane		352	C25H52	001560-78-7	45
3	Hexacosane		366	C26H54	000630-01-3	43
4	Undecane, 5-methyl-		170	C12H26	001632-70-8	41
5	Undecane, 2-methyl-		170	C12H26	007045-71-8	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

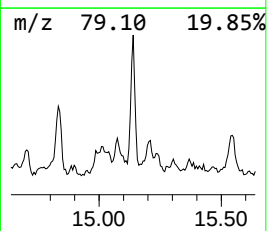
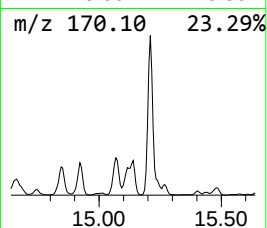
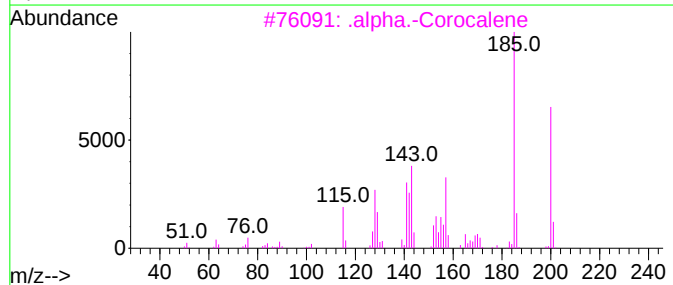
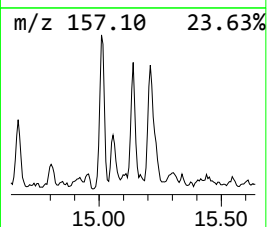
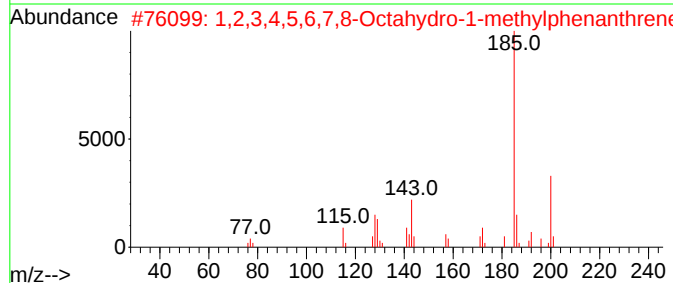
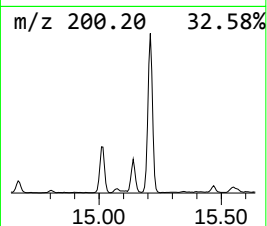
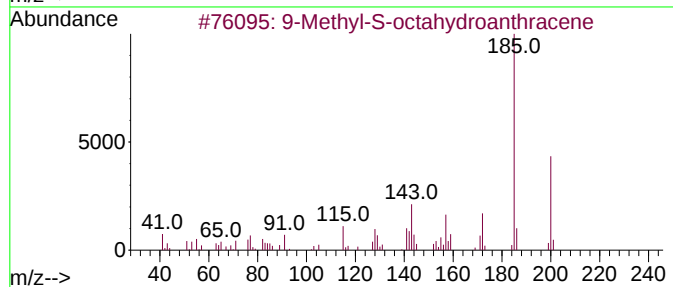
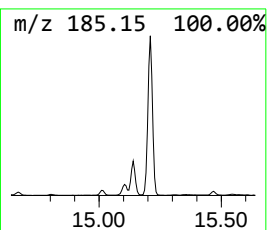
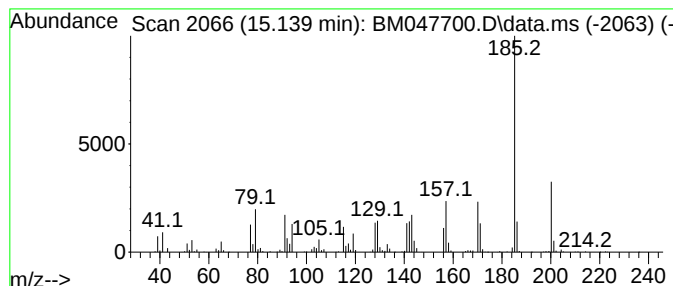
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 9-Methyl-S-octahydroanthracene Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.139	7.25 ng/ul	915895	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	76
2			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	70
3			.alpha.-Corocalene	200	C15H20	020129-39-9	64
4			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	58
5			Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
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Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

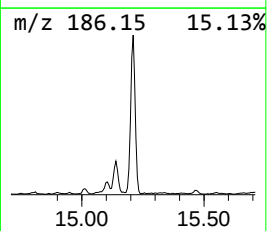
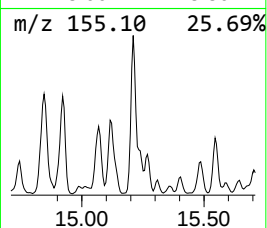
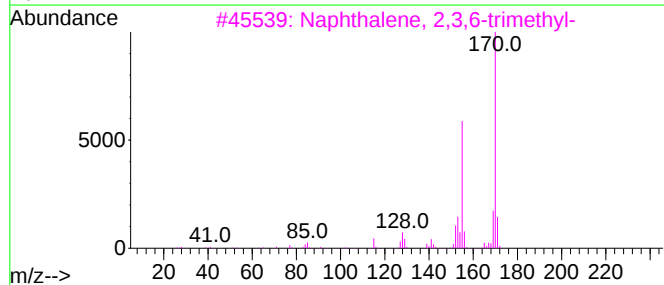
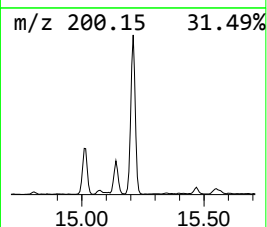
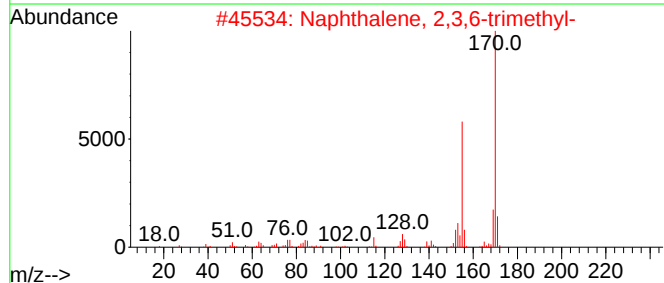
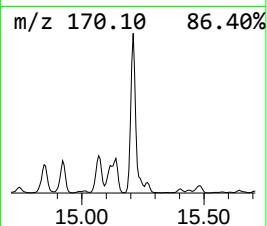
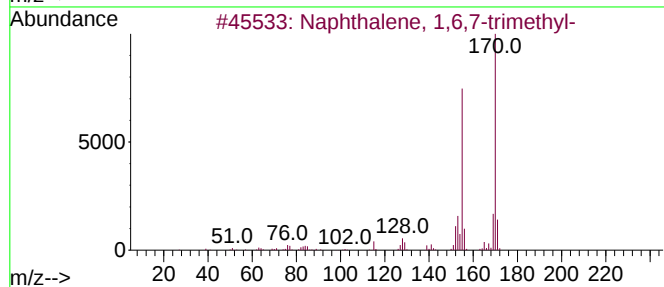
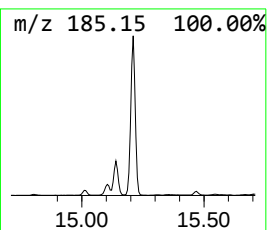
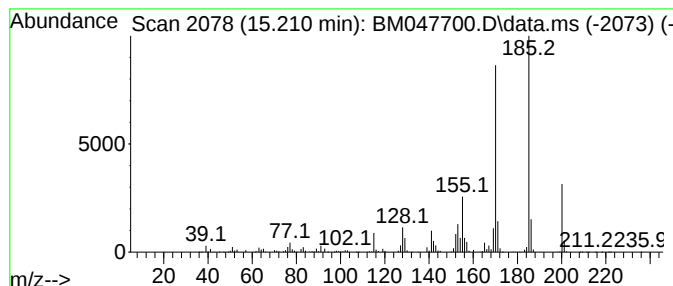
Instrument :
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ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 50 Naphthalene, 1,6,7-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.210	32.61 ng/ul	4118560	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	50
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
4	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

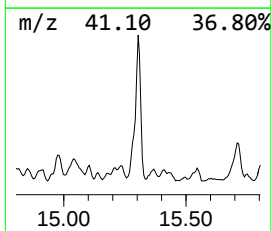
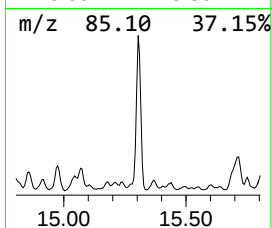
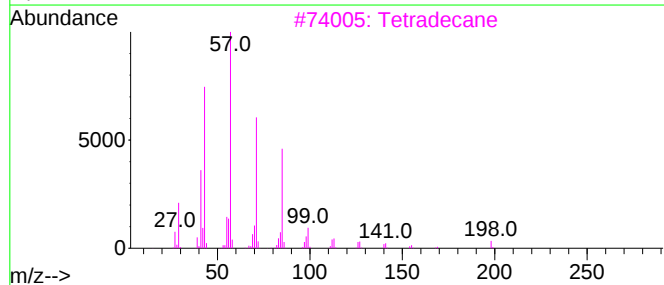
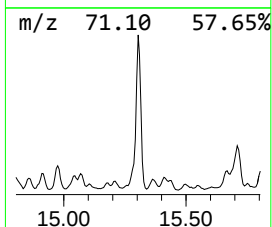
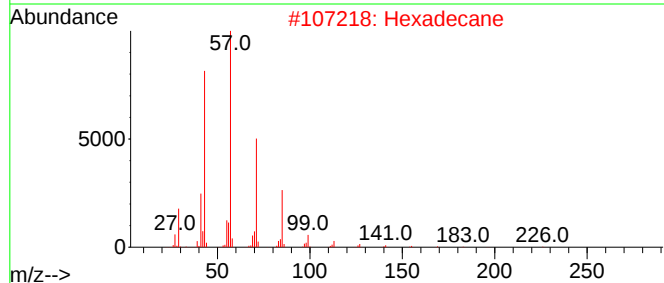
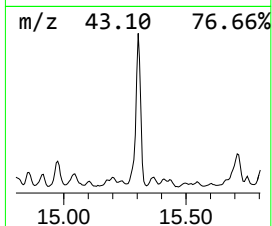
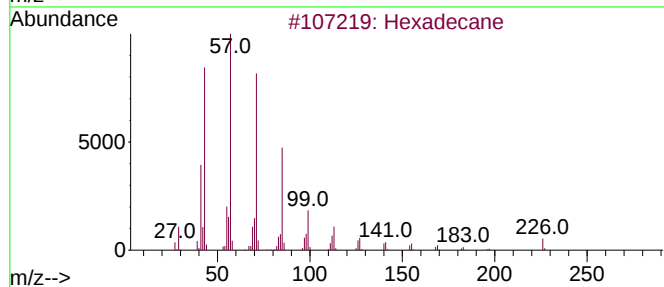
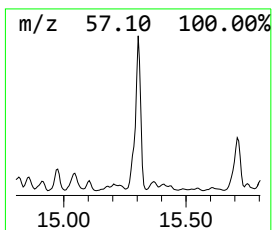
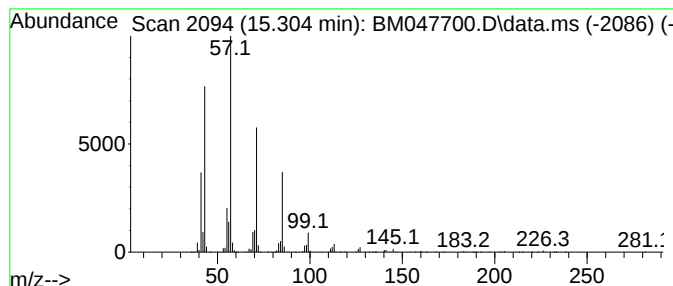
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Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.304	34.94 ng/ul	4412060	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	94
2			Hexadecane	226	C16H34	000544-76-3	91
3			Tetradecane	198	C14H30	000629-59-4	86
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

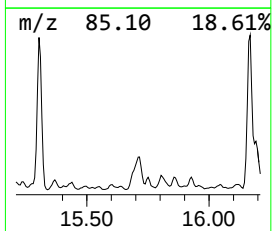
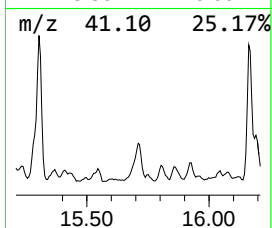
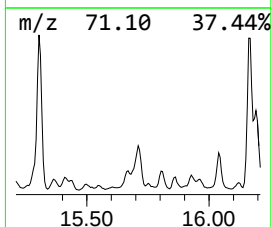
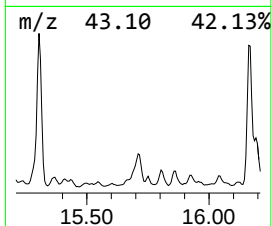
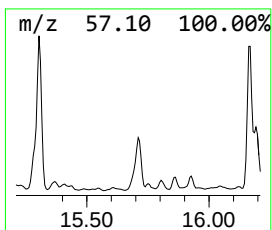
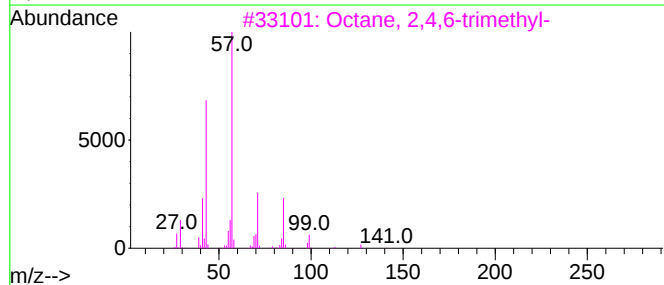
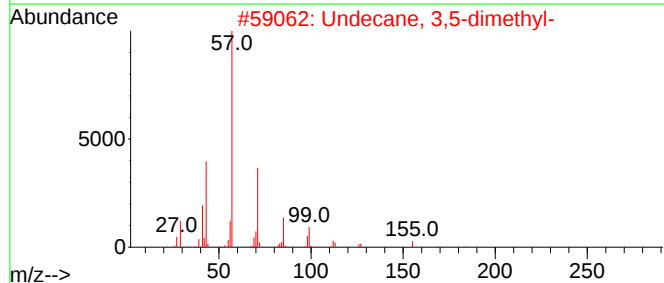
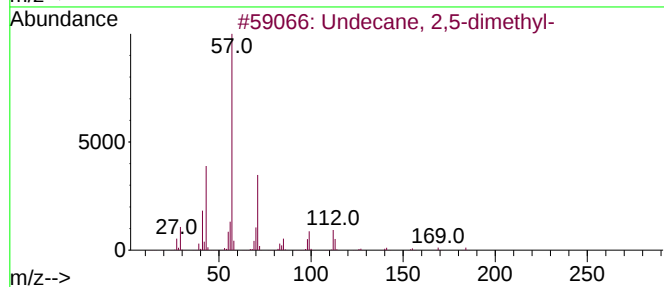
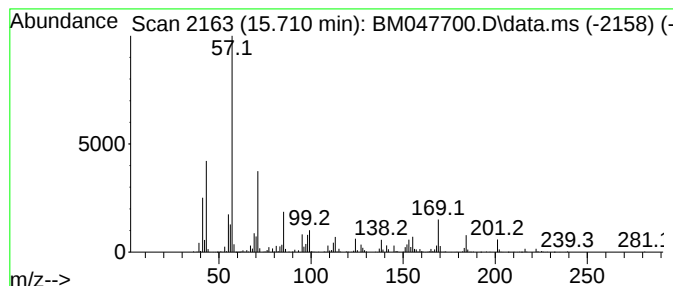
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.710	11.07 ng/ul	1397630	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 2,5-dimethyl-		184	C13H28	017301-22-3	58
2	Undecane, 3,5-dimethyl-		184	C13H28	017312-81-1	52
3	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	50
4	Dodecane, 3-methyl-		184	C13H28	017312-57-1	49
5	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

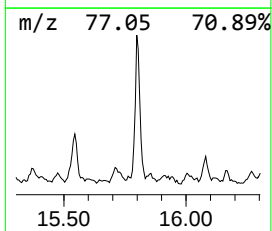
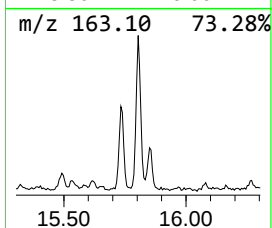
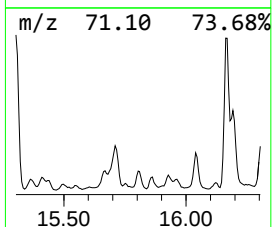
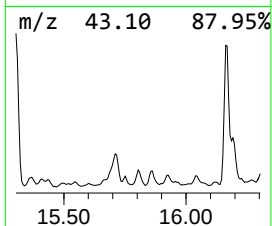
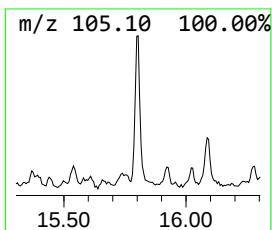
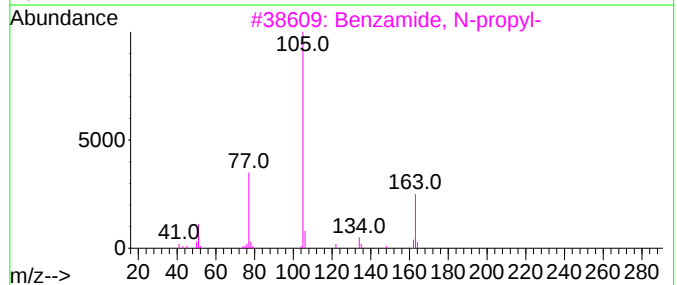
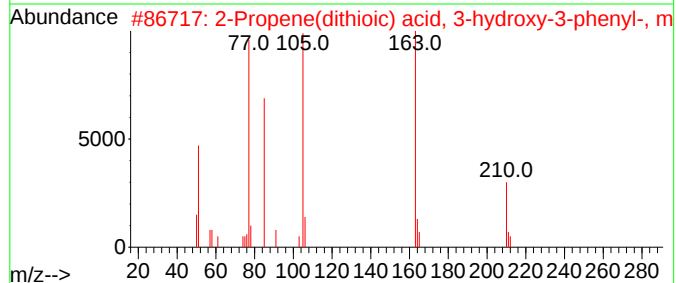
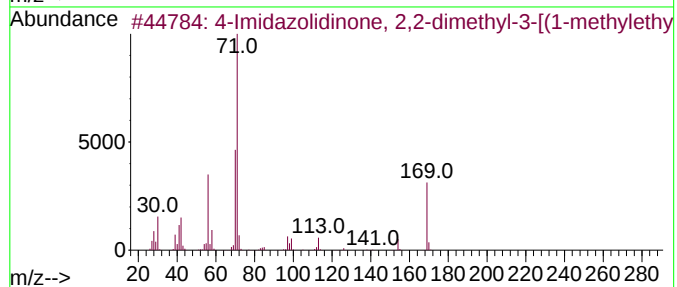
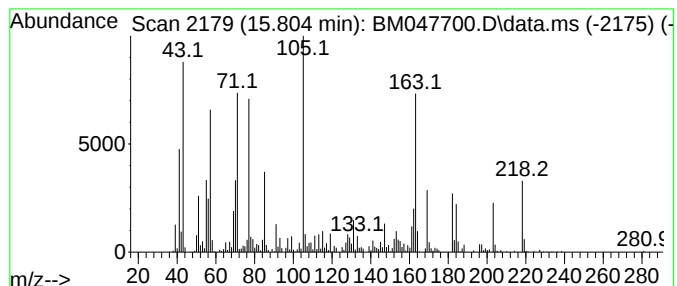
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 53 unknown-05 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.			
15.804	8.51 ng/ul	1074710	Acenaphthene-d10	14.445			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Imidazolidinone, 2,2-dimethyl-...	169	C8H15N3O	142071-78-1	35
2			2-Propene(dithioic) acid, 3-hydr...	210	C10H10S2	013636-68-5	30
3			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	25
4			4-Benzoyl-1-piperazinecarbaldehyde	218	C12H14N2O2	1000332-13-9	25
5			Acetamide, N-(3,4-dichlorophenyl...	219	C8H7Cl2NO2	086412-49-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

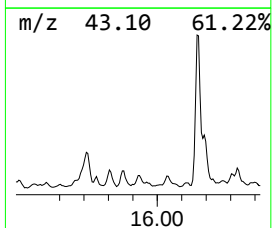
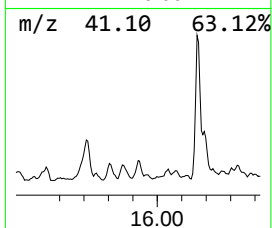
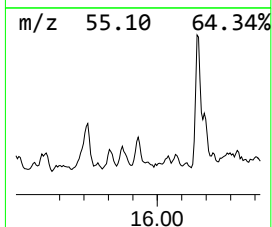
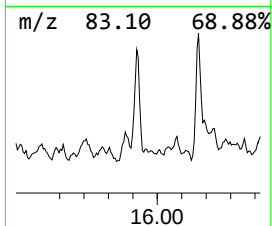
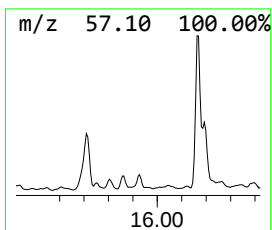
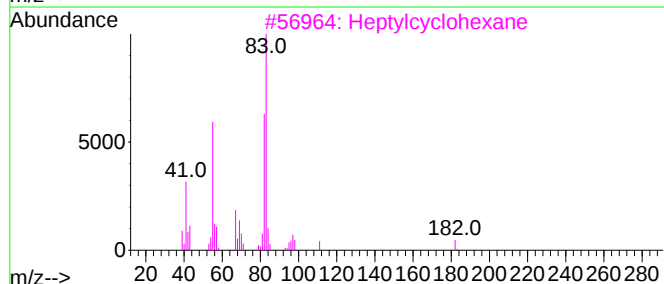
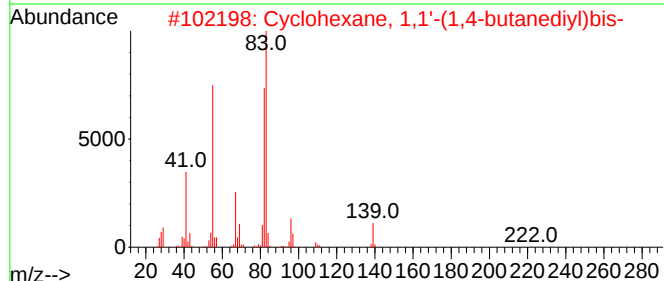
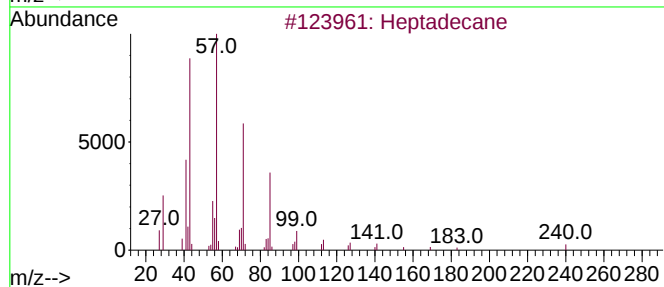
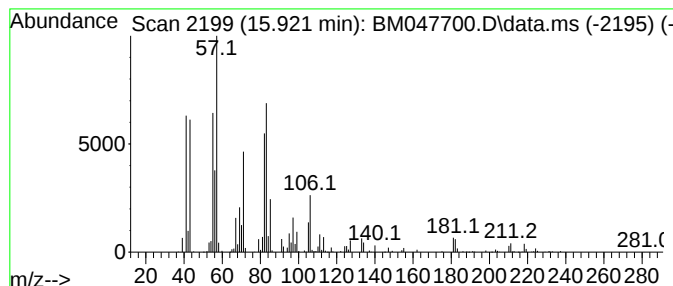
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.922	6.41 ng/ul	884898	Phenanthrene-d10			17.186
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	38
2	Cyclohexane, 1,1'-(1,4-butanediyl)bis-		222	C16H30	006165-44-2	38
3	Heptylcyclohexane		182	C13H26	005617-41-4	38
4	Silane, trichlorooctadecyl-		386	C18H37Cl3Si	000112-04-9	38
5	Methoxyacetic acid, 2-tetradecyl-		286	C17H34O3	1000282-04-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

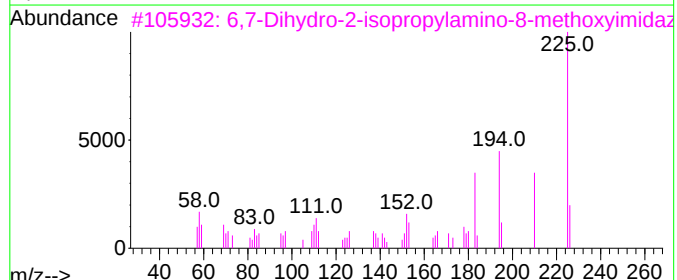
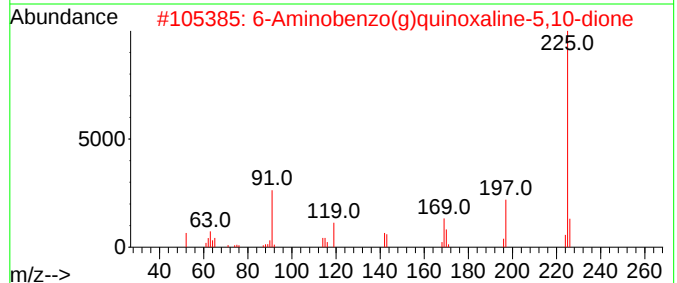
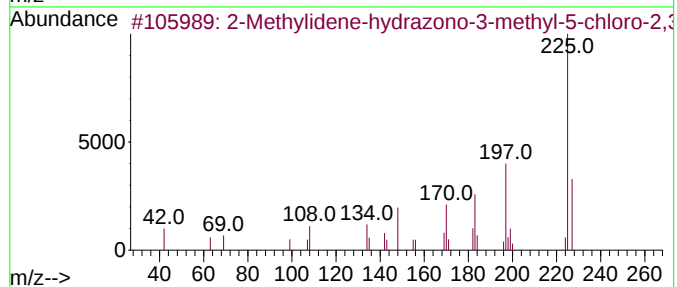
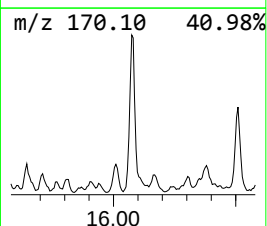
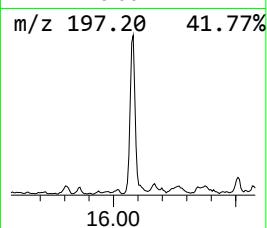
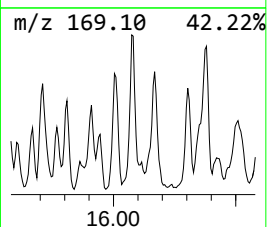
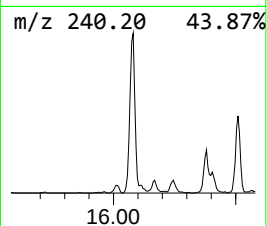
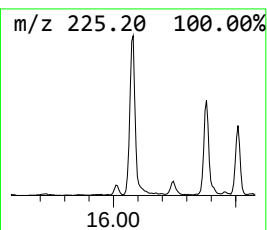
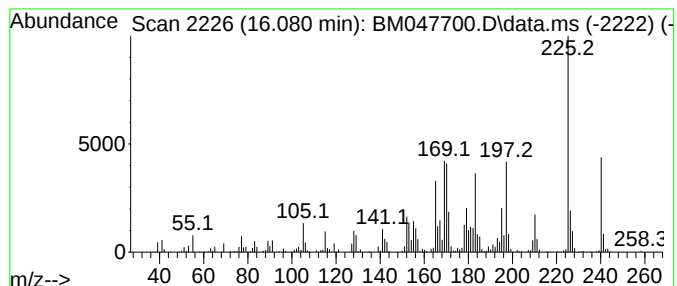
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Quant Title : SVOA CALIBRATION

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

Peak Number 55 unknown-06 Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.080	8.19 ng/ul	1131740	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	45
2			6-Aminobenzo(g)quinoxaline-5,10-...	225	C12H7N3O2	072225-24-2	43
3			6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	42
4			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	35
5			2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

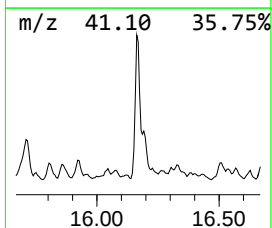
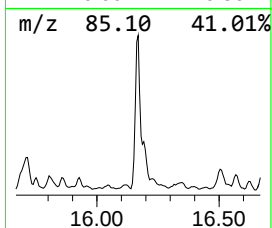
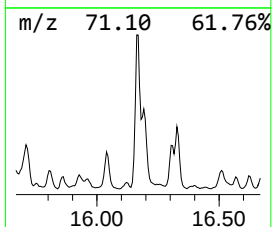
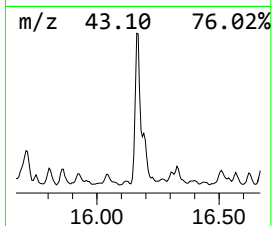
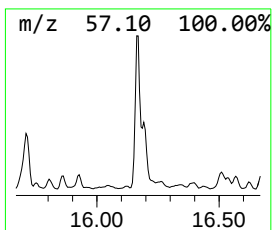
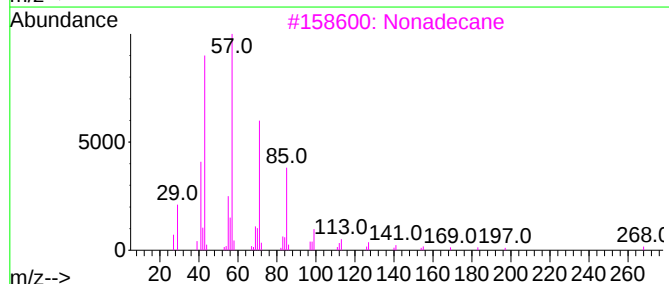
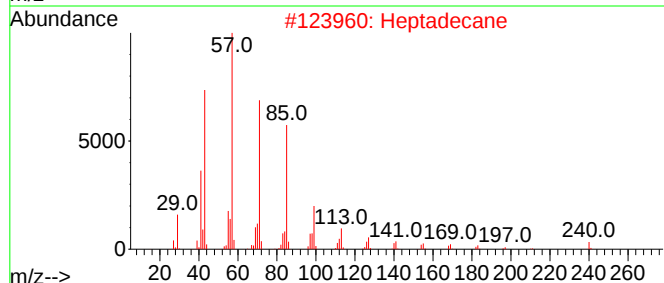
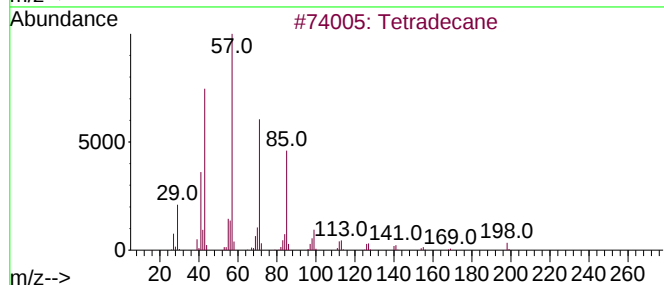
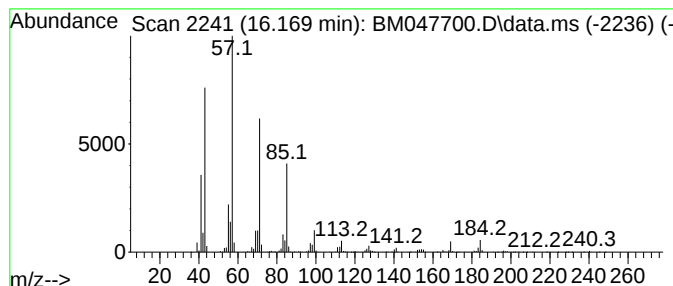
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.169	28.51 ng/ul	3938210	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	95
2		Heptadecane	240	C17H36	000629-78-7	94
3		Nonadecane	268	C19H40	000629-92-5	91
4		Heptadecane	240	C17H36	000629-78-7	87
5		Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

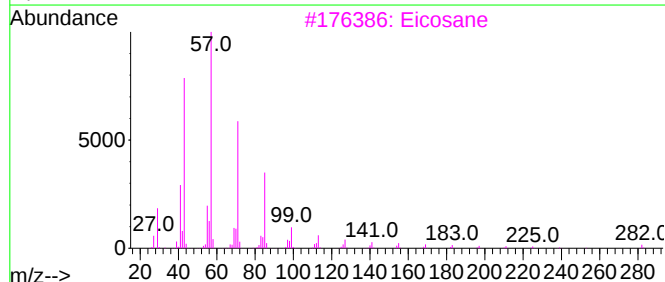
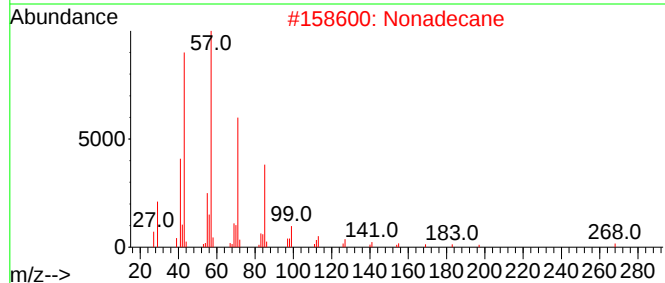
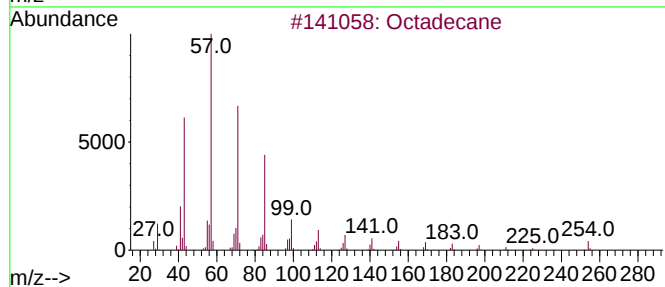
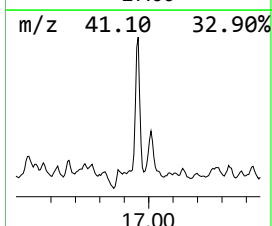
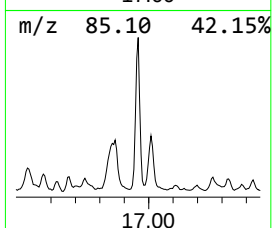
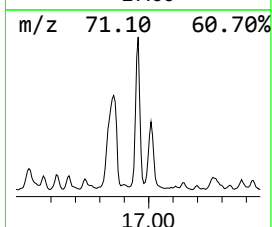
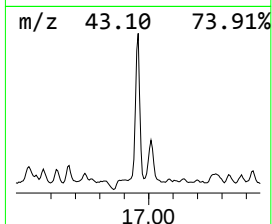
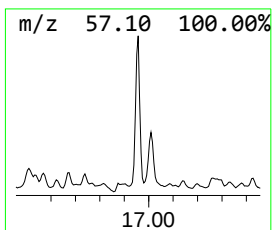
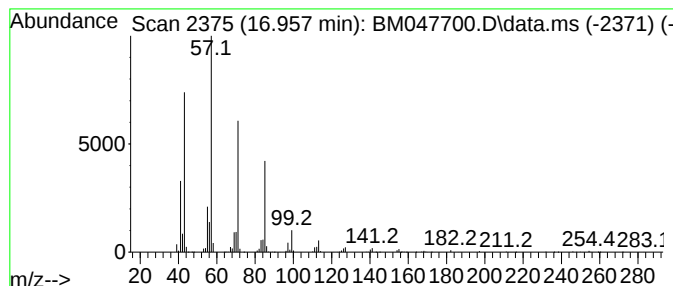
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.957	20.34 ng/ul	2809860	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	90
2			Nonadecane	268	C19H40	000629-92-5	90
3			Eicosane	282	C20H42	000112-95-8	86
4			Heptacosane	380	C27H56	000593-49-7	86
5			Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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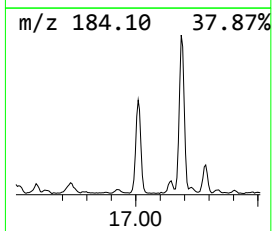
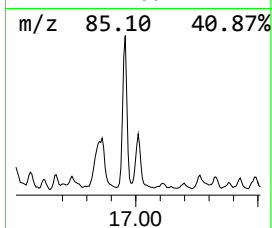
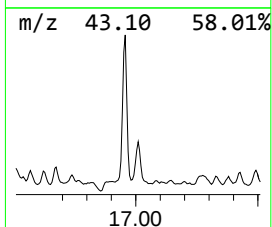
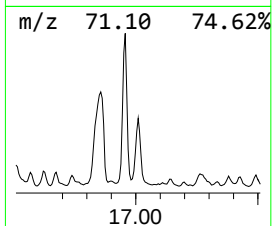
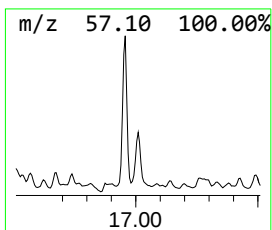
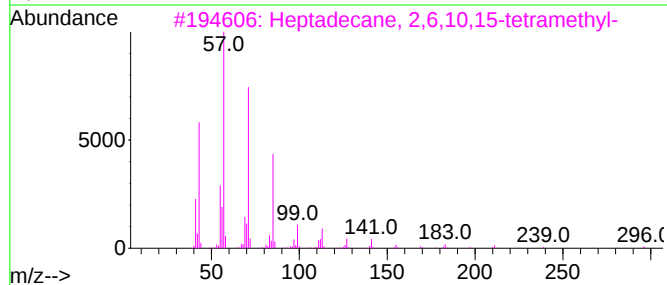
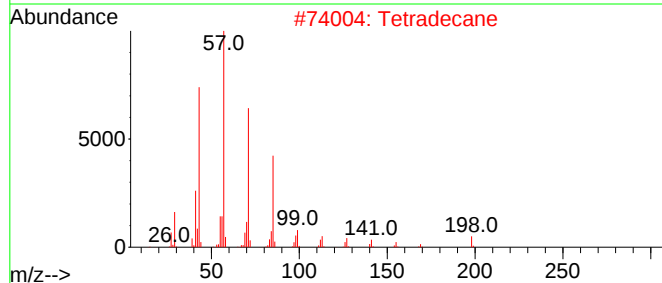
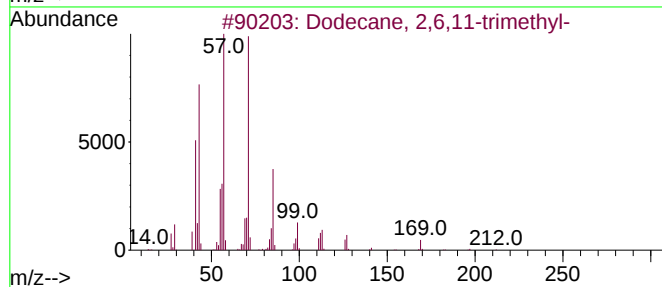
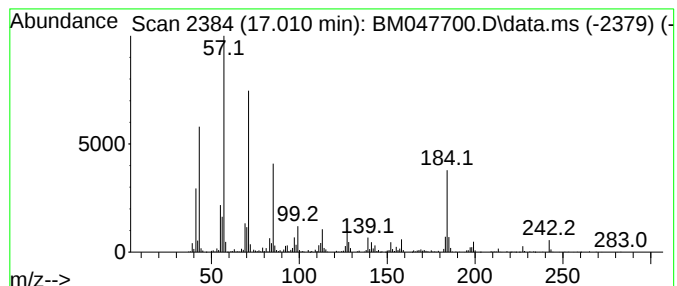
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.010	13.47 ng/ul	1861100	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	83
2		Tetradecane	198	C14H30	000629-59-4	81
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	74
4		Tetracosane	338	C24H50	000646-31-1	72
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

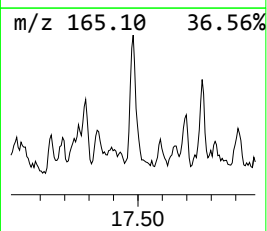
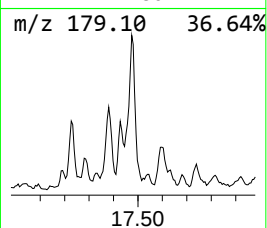
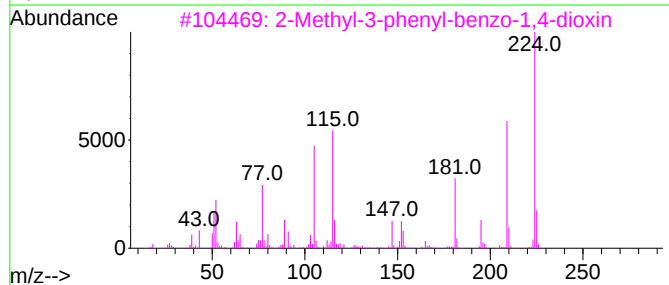
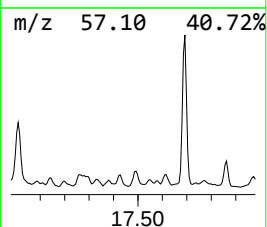
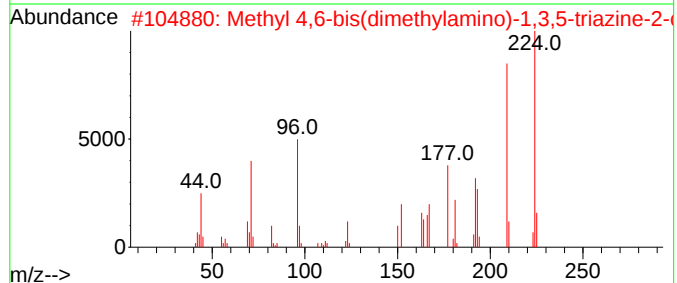
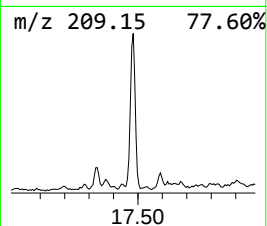
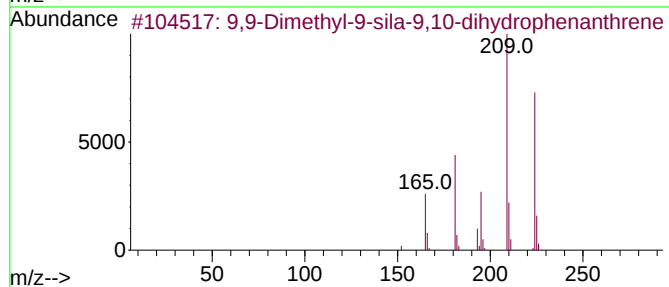
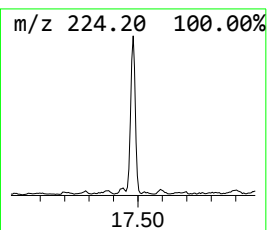
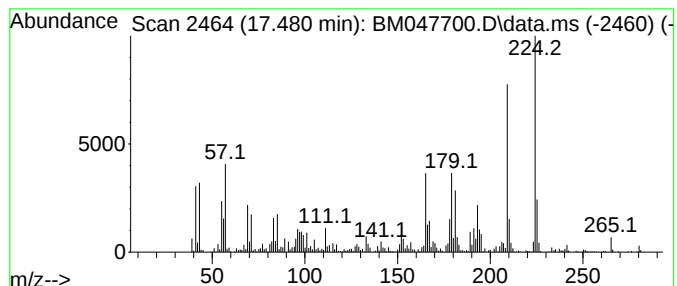
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 59 9,9-Dimethyl-9-sila-9,10-di... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.480	8.39 ng/ul	1158960	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si	187328-55-8	81
2	Methyl 4,6-bis(dimethylamino)-1,...	224	C9H16N6O	1000101-98-0	64
3	2-Methyl-3-phenyl-benzo-1,4-dioxin	224	C15H12O2	079792-91-9	53
4	7-Methyl-6,8-bis(methylthio)pyrr...	224	C10H12N2S2	084201-40-1	52
5	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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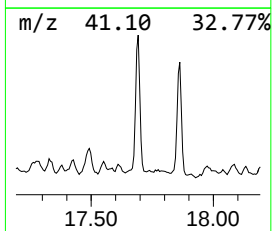
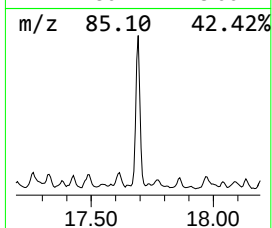
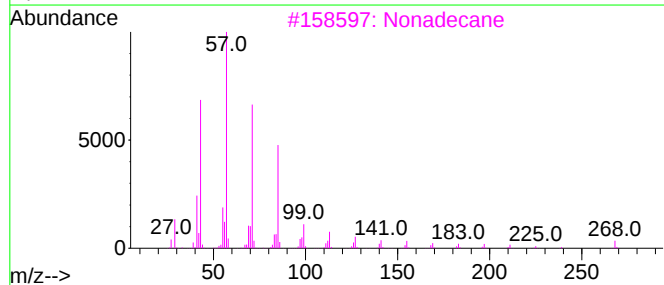
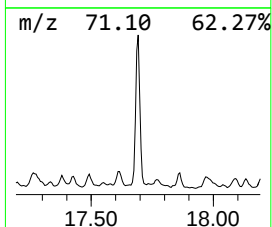
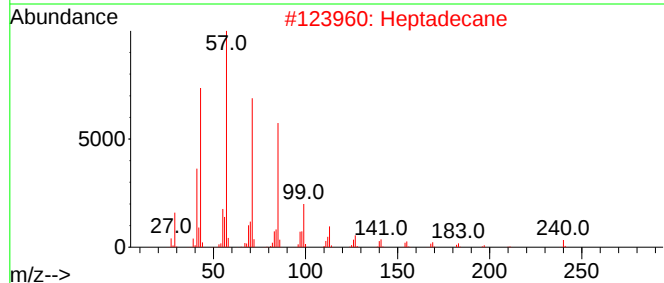
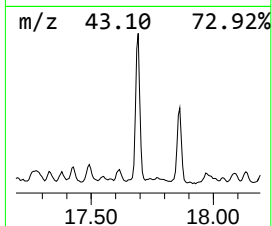
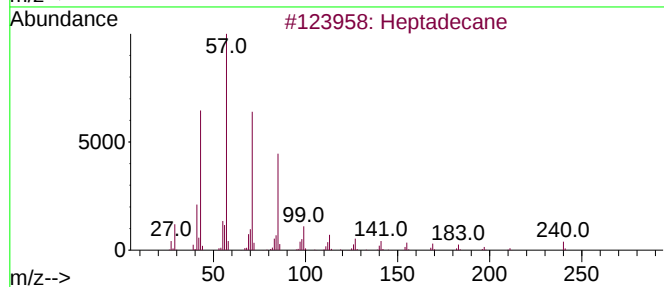
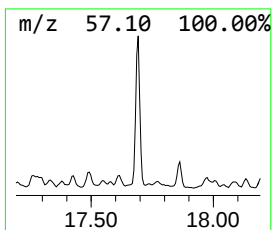
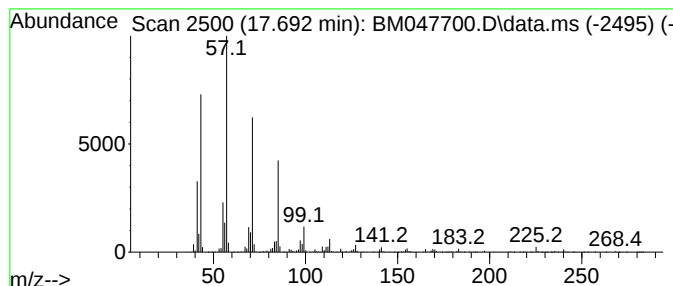
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.692	17.84 ng/ul	2463820	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	97
2		Heptadecane	240	C17H36	000629-78-7	96
3		Nonadecane	268	C19H40	000629-92-5	96
4		Octadecane	254	C18H38	000593-45-3	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

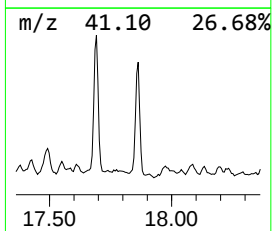
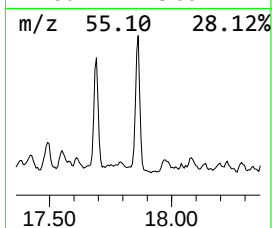
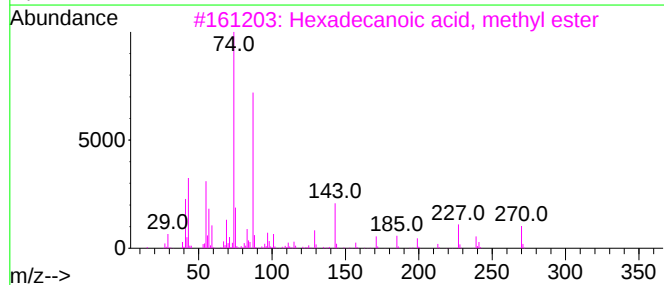
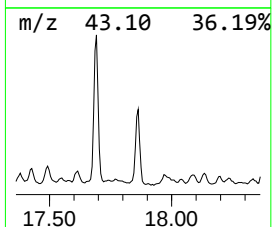
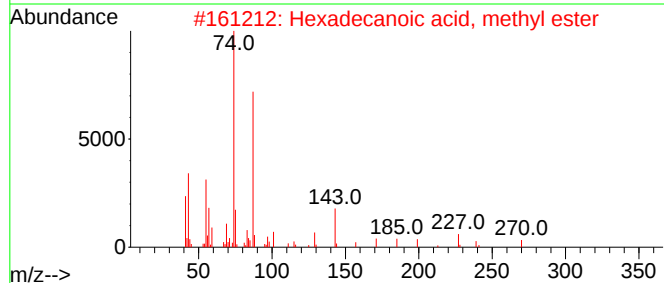
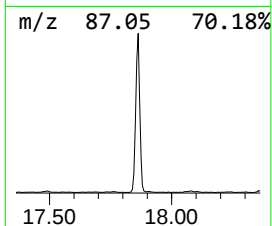
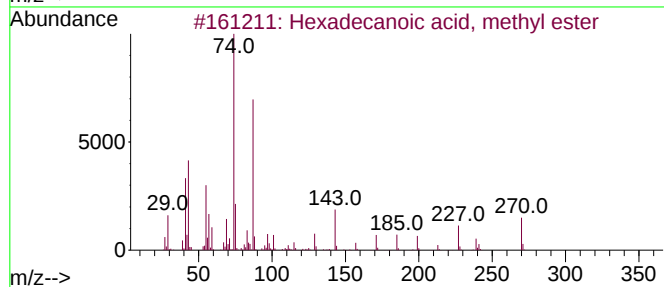
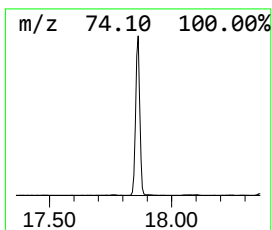
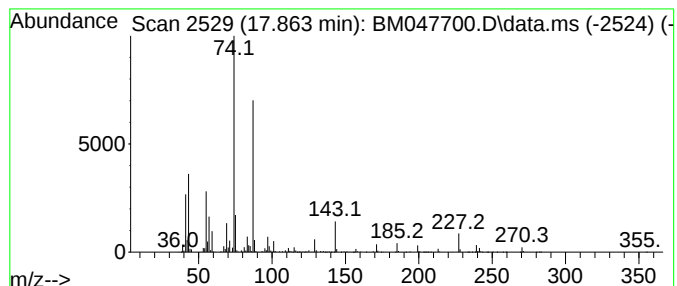
Instrument :
BNA_M
ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 61 Hexadecanoic acid, methyl e... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.863	16.15 ng/ul	2230380	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
3	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
5	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

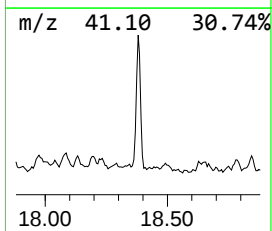
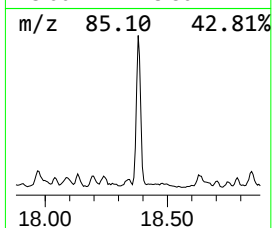
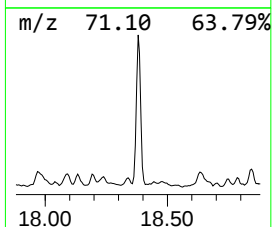
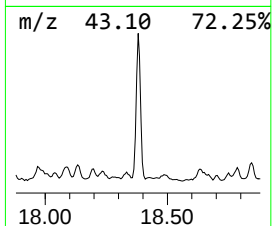
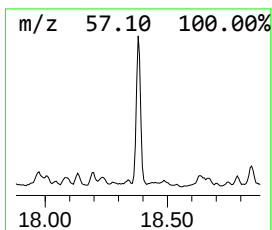
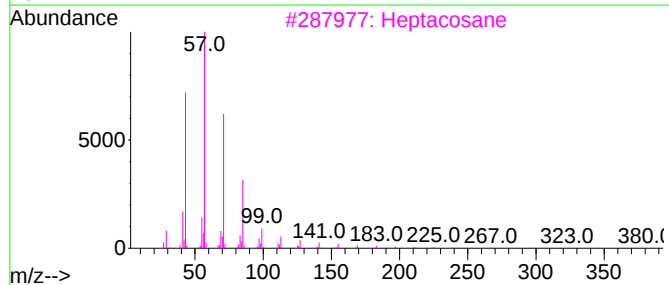
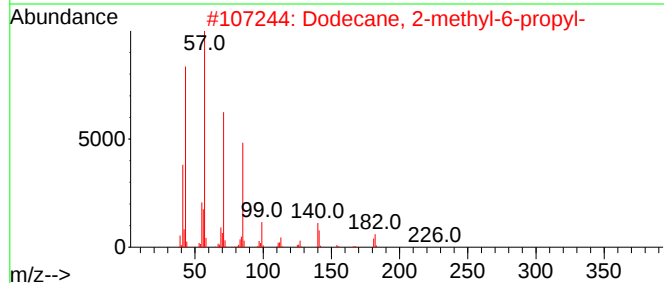
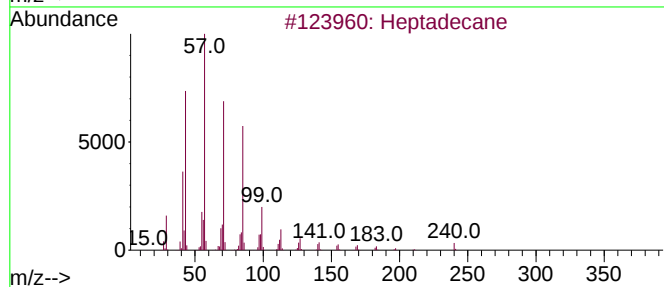
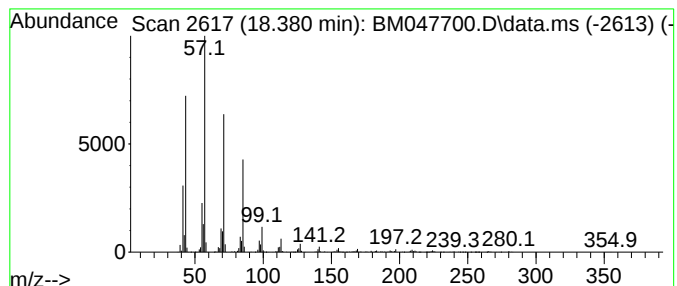
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BNA_M
ClientSampleId :
DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.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 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.380	13.19 ng/ul	1821600	Phenanthrene-d10		17.186	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	93
2	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	93
3	Heptacosane		380	C27H56	000593-49-7	91
4	Nonadecane		268	C19H40	000629-92-5	91
5	Hexadecane		226	C16H34	000544-76-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

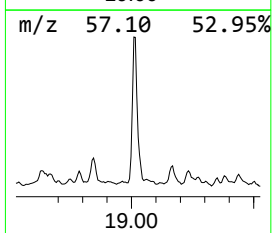
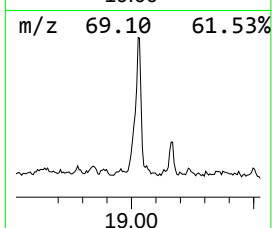
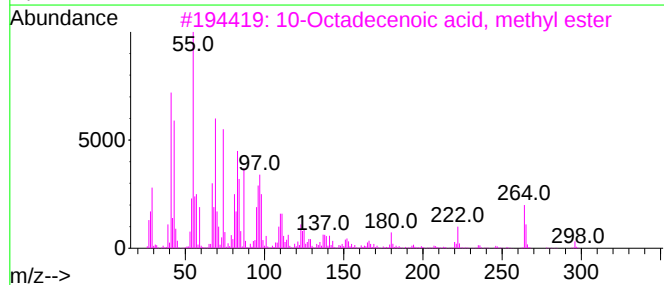
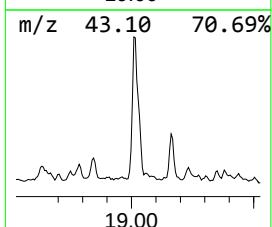
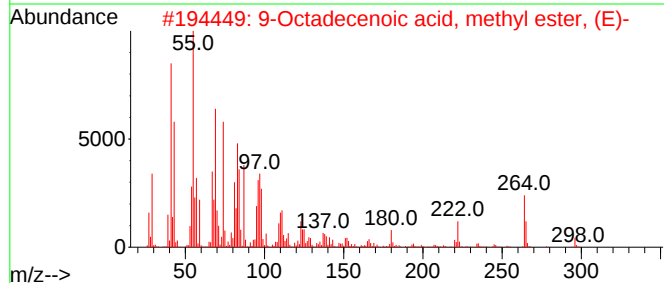
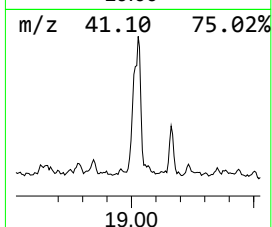
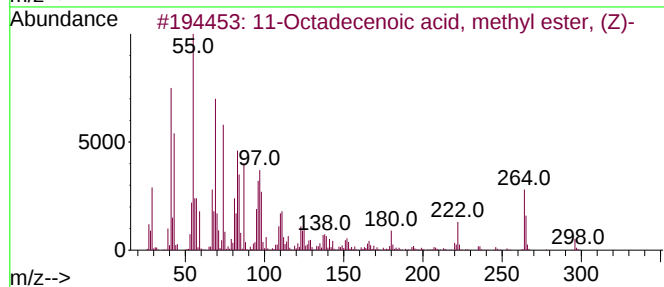
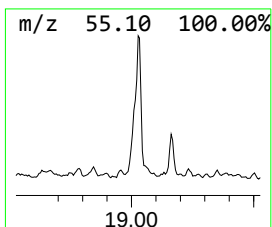
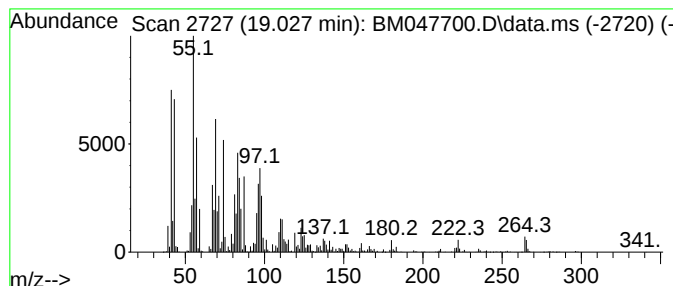
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 11-Octadecenoic acid, methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.027	23.29 ng/ul	3216840	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
2			9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
3			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	98
5			6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

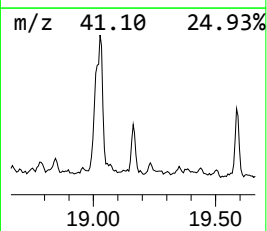
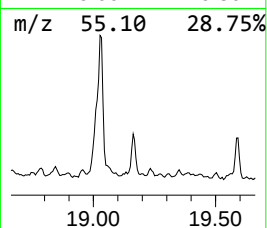
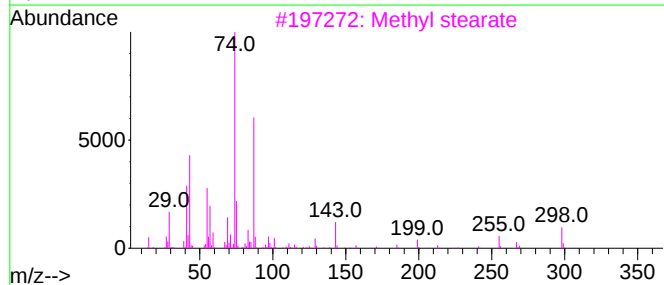
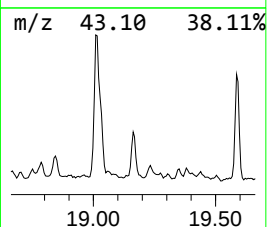
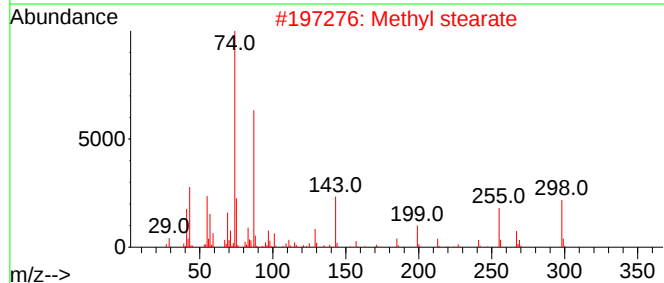
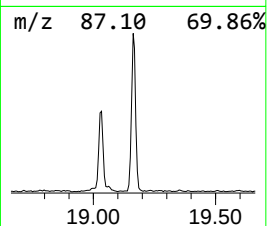
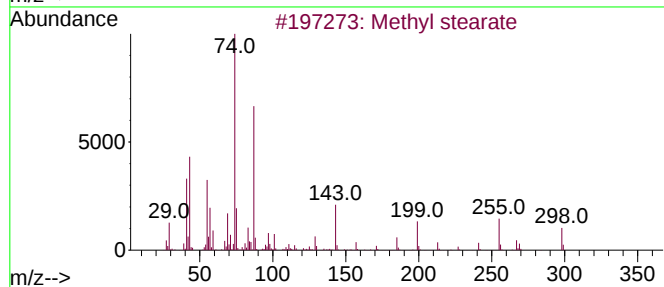
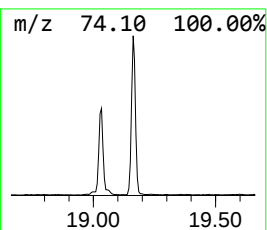
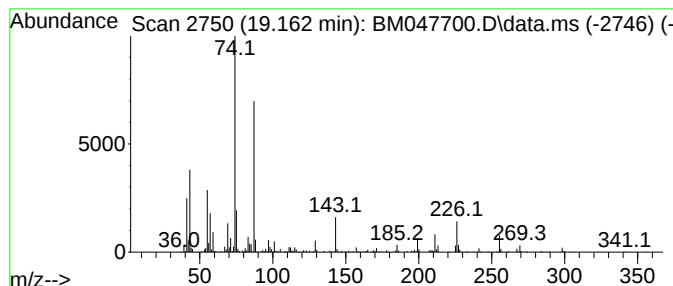
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Quant Title : SVOA CALIBRATION

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

Peak Number 64 Methyl stearate Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.162	7.00 ng/ul	966321	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Methyl stearate	298	C19H38O2	000112-61-8	97
3			Methyl stearate	298	C19H38O2	000112-61-8	96
4			Methyl stearate	298	C19H38O2	000112-61-8	93
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

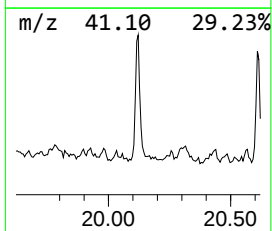
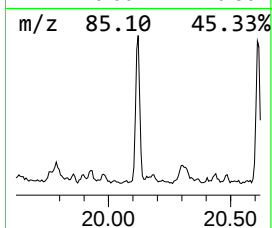
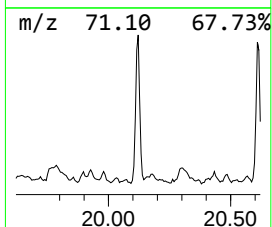
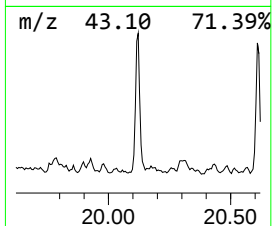
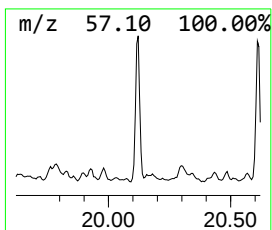
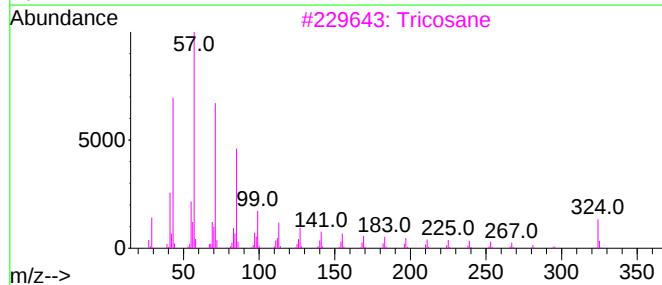
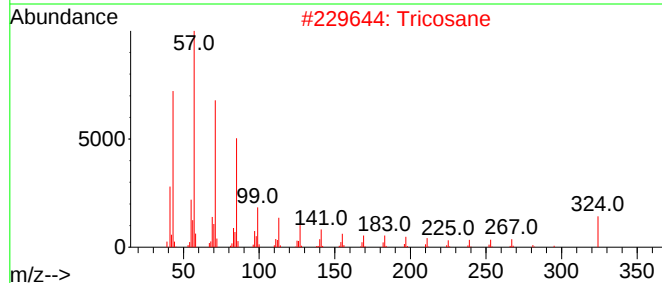
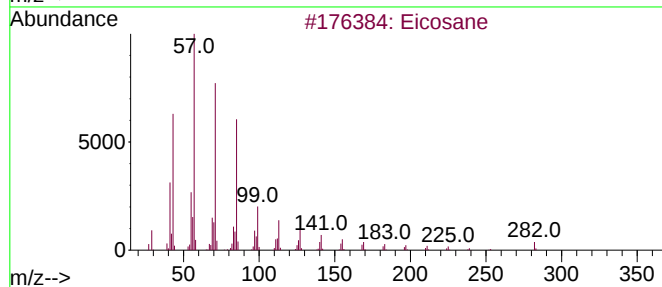
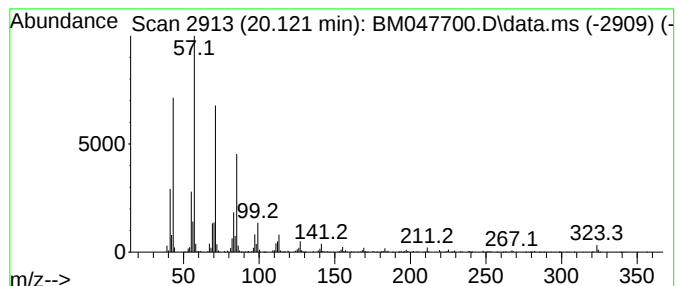
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.121	8.14 ng/ul	1150860	Chrysene-d12		21.403	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Tricosane		324	C23H48	000638-67-5	97
3	Tricosane		324	C23H48	000638-67-5	94
4	Nonadecane		268	C19H40	000629-92-5	93
5	Pentadecane		212	C15H32	000629-62-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

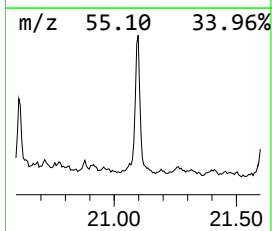
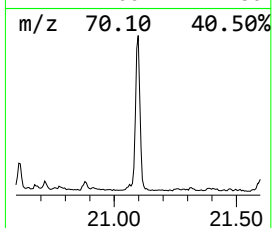
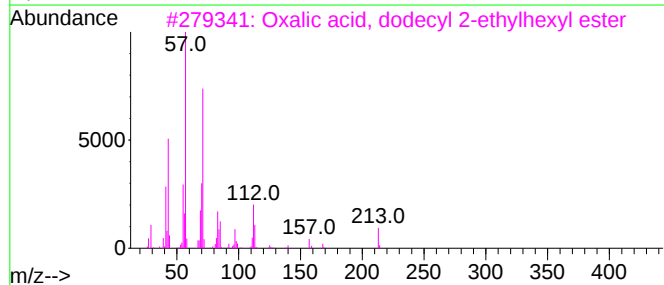
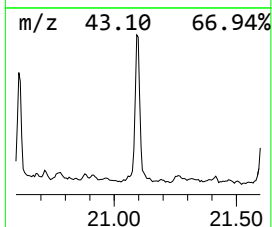
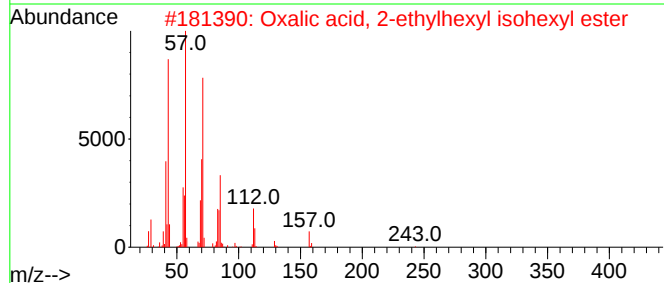
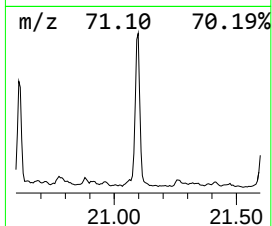
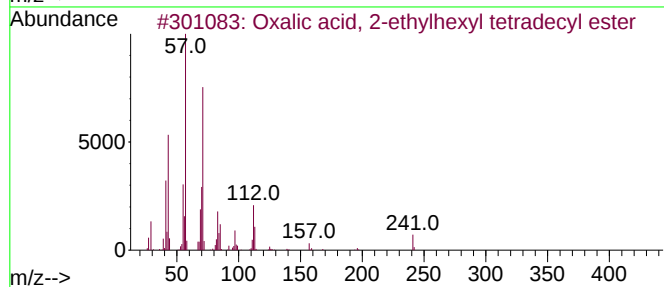
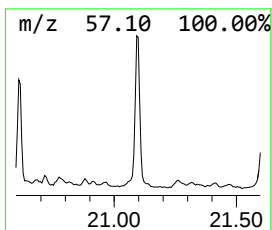
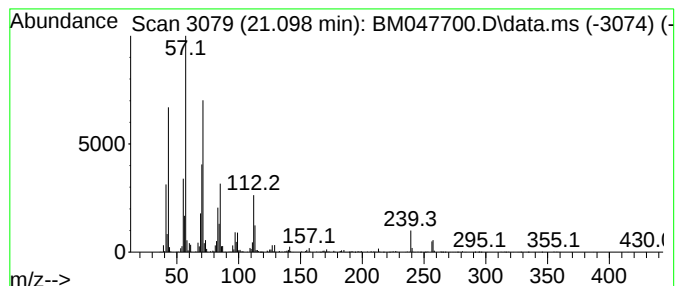
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Quant Title : SVOA CALIBRATION

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

Peak Number 66 Oxalic acid, 2-ethylhexyl t... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	12.90 ng/ul	1823570	Chrysene-d12	21.403

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	80
2			Oxalic acid, 2-ethylhexyl isohex...	286	C16H30O4	1000309-38-8	64
3			Oxalic acid, dodecyl 2-ethylhexy...	370	C22H42O4	1000309-39-5	52
4			Decane, 2,3,4-trimethyl-	184	C13H28	062238-15-7	49
5			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
Data File : BM047700.D
Acq On : 28 Sep 2024 20:38
Operator : RC/JU
Sample : P3927-08DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDC93DL

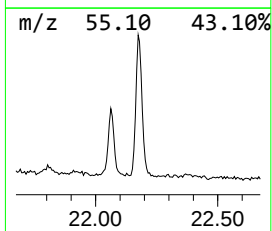
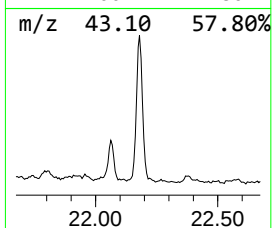
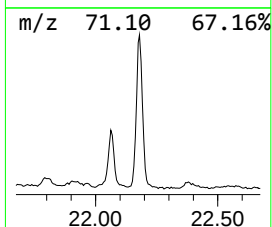
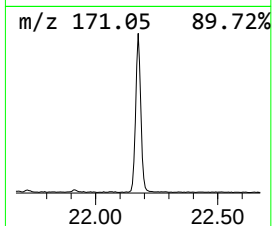
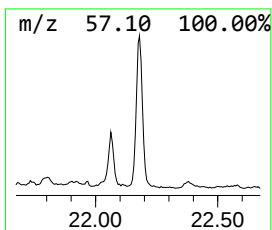
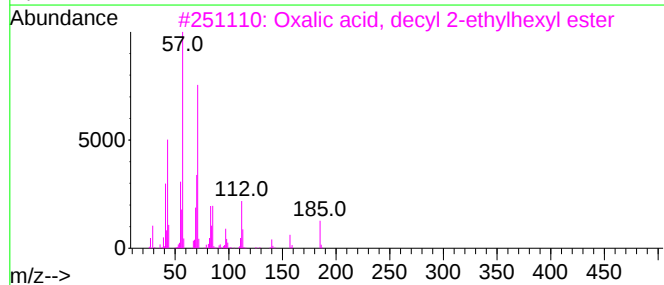
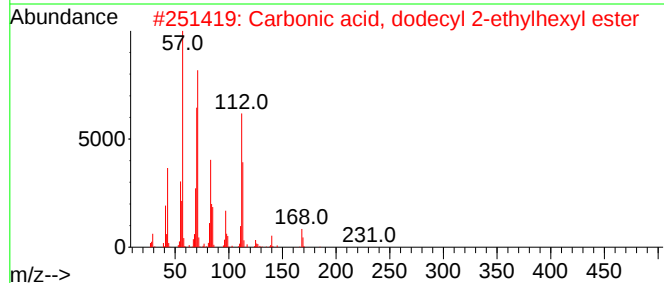
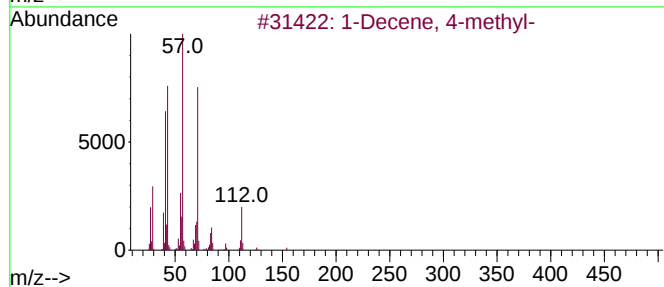
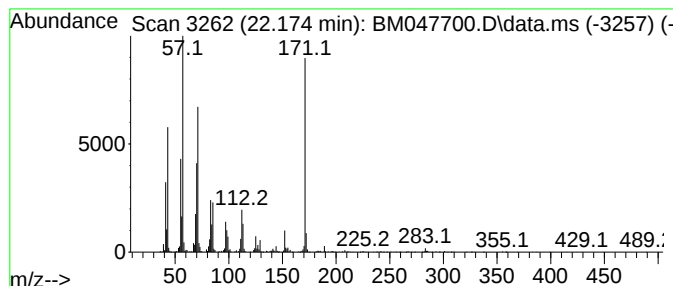
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Quant Title : SVOA CALIBRATION

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

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	13.59 ng/ul	1920890	Chrysene-d12	21.403

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decene, 4-methyl-	154	C11H22	013151-29-6	60
2			Carbonic acid, dodecyl 2-ethylhe...	342	C21H42O3	1000383-13-9	49
3			Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	49
4			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	38
5			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092824\
 Data File : BM047700.D
 Acq On : 28 Sep 2024 20:38
 Operator : RC/JU
 Sample : P3927-08DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDC93DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.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
(DEL) Alkane: S...	5.840	19.7	ng/ul	3472820	1	7.816	3521700	20.0
(DEL) Alkane: S...	6.381	6.2	ng/ul	1094160	1	7.816	3521700	20.0
(DEL) Alkane: C...	6.463	7.7	ng/ul	1348440	1	7.816	3521700	20.0
(DEL) Alkane: S...	6.816	15.9	ng/ul	2800950	1	7.816	3521700	20.0
(DEL) Alkane: S...	6.875	11.8	ng/ul	2083160	1	7.816	3521700	20.0
Benzene, 1-ethy...	6.910	12.5	ng/ul	2201400	1	7.816	3521700	20.0
Mesitylene	7.046	8.0	ng/ul	1400840	1	7.816	3521700	20.0
Benzene, 1,2,4-...	7.210	8.3	ng/ul	1465580	1	7.816	3521700	20.0
2-Heptanone, 4,...	7.246	8.3	ng/ul	1459720	1	7.816	3521700	20.0
(DEL) Alkane: S...	7.451	79.5	ng/ul	14002300	1	7.816	3521700	20.0
Carbonic acid, ...	7.893	9.0	ng/ul	1592050	1	7.816	3521700	20.0
Benzene, 1,2,3-...	7.934	9.1	ng/ul	1608890	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.016	6.2	ng/ul	1083200	1	7.816	3521700	20.0
unknown-01	8.040	6.5	ng/ul	1145930	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.087	5.1	ng/ul	905266	1	7.816	3521700	20.0
(DEL) Alkane: C...	8.116	6.2	ng/ul	1095230	1	7.816	3521700	20.0
Cyclohexanone, ...	8.245	13.1	ng/ul	2308600	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.298	8.8	ng/ul	1544560	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.357	17.3	ng/ul	3046930	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.416	8.6	ng/ul	1516470	1	7.816	3521700	20.0
Benzene, 1,4-di...	8.457	11.4	ng/ul	2013230	1	7.816	3521700	20.0
(DEL) Alkane: S...	8.593	14.4	ng/ul	2541780	1	7.816	3521700	20.0
Benzene, 1-meth...	8.622	7.2	ng/ul	1261080	1	7.816	3521700	20.0
o-Cymene	8.816	10.1	ng/ul	1774830	1	7.816	3521700	20.0
(DEL) Alkane: S...	9.051	63.9	ng/ul	11249500	1	7.816	3521700	20.0
Benzene, 1-meth...	9.175	12.1	ng/ul	2133140	1	7.816	3521700	20.0
(DEL) Alkane: S...	9.898	2.4	ng/ul	1907730	2	10.610	15720900	20.0
(DEL) Alkane: S...	11.122	2.1	ng/ul	1672730	2	10.610	15720900	20.0
(DEL) Alkane: S...	11.645	2.5	ng/ul	2000270	2	10.610	15720900	20.0
(DEL) Alkane: S...	12.039	4.5	ng/ul	3577930	2	10.610	15720900	20.0
(DEL) Alkane: S...	12.686	14.3	ng/ul	1800050	3	14.445	2525640	20.0
1-Iodo-2-methyl...	12.857	6.8	ng/ul	858165	3	14.445	2525640	20.0
unknown-02	13.504	10.8	ng/ul	1362300	3	14.445	2525640	20.0
Naphthalene, 1,...	13.633	15.8	ng/ul	1992370	3	14.445	2525640	20.0
Butanoic acid, ...	13.716	8.0	ng/ul	1015040	3	14.445	2525640	20.0
Naphthalene, 1,...	13.769	12.3	ng/ul	1549120	3	14.445	2525640	20.0
(DEL) Alkane: S...	13.945	8.2	ng/ul	1040840	3	14.445	2525640	20.0
unknown-03	14.086	13.9	ng/ul	1761470	3	14.445	2525640	20.0
(DEL) Alkane: S...	14.363	104.3	ng/ul	13171600	3	14.445	2525640	20.0
unknown-04	14.527	8.7	ng/ul	1101760	3	14.445	2525640	20.0
Naphthalene, 2,...	14.845	7.8	ng/ul	982753	3	14.445	2525640	20.0
Naphthalene, 1,...	14.916	7.0	ng/ul	878188	3	14.445	2525640	20.0
(DEL) Alkane: S...	14.980	7.5	ng/ul	942715	3	14.445	2525640	20.0
9-Methyl-S-octa...	15.139	7.3	ng/ul	915895	3	14.445	2525640	20.0
Naphthalene, 1,...	15.210	32.6	ng/ul	4118560	3	14.445	2525640	20.0
(DEL) Alkane: S...	15.304	34.9	ng/ul	4412060	3	14.445	2525640	20.0
(DEL) Alkane: S...	15.710	11.1	ng/ul	1397630	3	14.445	2525640	20.0
unknown-05	15.804	8.5	ng/ul	1074710	3	14.445	2525640	20.0
(DEL) Alkane: S...	15.922	6.4	ng/ul	884898	4	17.186	2762840	20.0
unknown-06	16.080	8.2	ng/ul	1131740	4	17.186	2762840	20.0
(DEL) Alkane: S...	16.169	28.5	ng/ul	3938210	4	17.186	2762840	20.0
(DEL) Alkane: S...	16.957	20.3	ng/ul	2809860	4	17.186	2762840	20.0

(DEL) Alkane: S...	17.010	13.5	ng/ul	1861100	4	17.186	2762840	20.0
9,9-Dimethyl-9-...	17.480	8.4	ng/ul	1158960	4	17.186	2762840	20.0
(DEL) Alkane: S...	17.692	17.8	ng/ul	2463820	4	17.186	2762840	20.0
Hexadecanoic ac...	17.863	16.1	ng/ul	2230380	4	17.186	2762840	20.0
(DEL) Alkane: S...	18.380	13.2	ng/ul	1821600	4	17.186	2762840	20.0
11-Octadecenoic...	19.027	23.3	ng/ul	3216840	4	17.186	2762840	20.0
Methyl stearate	19.162	7.0	ng/ul	966321	4	17.186	2762840	20.0
(DEL) Alkane: S...	20.121	8.1	ng/ul	1150860	5	21.403	2827290	20.0
Oxalic acid, 2-...	21.098	12.9	ng/ul	1823570	5	21.403	2827290	20.0
(DEL) Alkane: S...	22.174	13.6	ng/ul	1920890	5	21.403	2827290	20.0

Instrument :

BNA_M

ClientSampleId :

DDC93DL