

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047009.D
 Acq On : 06 Aug 2024 02:47
 Operator : RC/JU
 Sample : P3171-05DL 30X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVW9DL

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-BM080224.MA.M

Title : SVOA CALIBRATION

Signal : TIC: BM047009.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.463	431	438	446	rVB2	265000	377853	2.78%	0.720%
2	5.775	486	491	495	rBV	188627	264976	1.95%	0.505%
3	5.987	522	527	532	rBV	76236	109025	0.80%	0.208%
4	6.063	532	540	543	rVV3	59476	133588	0.98%	0.255%
5	6.416	597	600	604	rVV	109562	147650	1.09%	0.281%
6	6.469	604	609	612	rVV	126384	187087	1.38%	0.357%
7	6.504	612	615	619	rVV	127084	184651	1.36%	0.352%
8	6.575	619	627	634	rVV4	168401	407718	3.00%	0.777%
9	6.634	634	637	642	rVV	102488	148599	1.09%	0.283%
10	6.793	660	664	669	rVV2	110888	178834	1.31%	0.341%
11	6.846	669	673	678	rVV3	78523	142251	1.05%	0.271%
12	6.940	684	689	700	rVV5	75202	172271	1.27%	0.328%
13	7.040	700	706	716	rVV2	752330	1242900	9.14%	2.369%
14	7.393	758	766	772	rBV2	775980	1207708	8.88%	2.302%
15	7.475	776	780	784	rBV	235251	353087	2.60%	0.673%
16	7.628	802	806	812	rVV	379613	560356	4.12%	1.068%
17	7.828	835	840	845	rBV2	78952	130017	0.96%	0.248%
18	7.881	845	849	853	rVV3	67470	116005	0.85%	0.221%
19	7.934	853	858	863	rVB3	157427	268914	1.98%	0.513%
20	7.993	863	868	871	rBV	93959	126638	0.93%	0.241%
21	8.028	871	874	877	rVV2	137813	213831	1.57%	0.408%
22	8.063	877	880	886	rVV	143524	227787	1.67%	0.434%
23	8.163	893	897	900	rVV2	121637	194971	1.43%	0.372%
24	8.193	900	902	909	rVB2	90526	147023	1.08%	0.280%
25	8.393	931	936	943	rVB	79902	125303	0.92%	0.239%
26	8.487	943	952	956	rBV2	110888	234229	1.72%	0.446%
27	8.622	970	975	983	rVB	689426	1003785	7.38%	1.913%
28	8.716	983	991	1001	rBV4	187766	503381	3.70%	0.960%
29	8.822	1001	1009	1013	rBV3	104512	199188	1.46%	0.380%
30	9.181	1063	1070	1077	rVB2	62605	110958	0.82%	0.212%
31	9.251	1077	1082	1087	rBV6	72906	171697	1.26%	0.327%
32	9.328	1087	1095	1098	rVV4	70117	179844	1.32%	0.343%
33	9.357	1098	1100	1109	rVB3	77136	158742	1.17%	0.303%
34	9.469	1109	1119	1124	rVB	82060	197373	1.45%	0.376%
35	9.716	1152	1161	1165	rBV3	70356	141814	1.04%	0.270%
36	9.869	1181	1187	1192	rBV2	73428	116016	0.85%	0.221%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Title : SVOA CALIBRATION

37	10.151	1225	1235	1241	rBV	1255486	2201459	16.19%	4.196%
38	10.292	1253	1259	1264	rBV3	200272	344935	2.54%	0.658%
39	10.551	1297	1303	1311	rBV	296128	483162	3.55%	0.921%
40	10.687	1315	1326	1332	rBV3	77969	182190	1.34%	0.347%
41	11.204	1408	1414	1419	rBV3	140506	271738	2.00%	0.518%
42	11.269	1419	1425	1432	rVB	294217	483654	3.56%	0.922%
43	11.445	1446	1455	1461	rBV2	239321	384599	2.83%	0.733%
44	11.610	1475	1483	1487	rBV	271500	482791	3.55%	0.920%
45	11.645	1487	1489	1494	rVV	189194	251227	1.85%	0.479%
46	11.786	1510	1513	1516	rVV3	74431	126699	0.93%	0.242%
47	11.828	1517	1520	1524	rVV	95111	157611	1.16%	0.300%
48	11.875	1524	1528	1537	rVB	335400	484182	3.56%	0.923%
49	12.051	1550	1558	1563	rBV3	64676	156334	1.15%	0.298%
50	12.281	1591	1597	1601	rBV4	75103	147510	1.08%	0.281%
51	12.457	1623	1627	1633	rVB2	88817	150630	1.11%	0.287%
52	12.598	1646	1651	1659	rVB2	85676	158668	1.17%	0.302%
53	12.810	1678	1687	1692	rBV2	89993	175288	1.29%	0.334%
54	12.892	1694	1701	1712	rBV	314731	496862	3.65%	0.947%
55	13.022	1719	1723	1727	rBV	77953	109772	0.81%	0.209%
56	13.210	1747	1755	1758	rBV6	69092	159613	1.17%	0.304%
57	13.333	1771	1776	1779	rBV	174535	251132	1.85%	0.479%
58	13.369	1779	1782	1786	rVV	85465	118098	0.87%	0.225%
59	13.422	1786	1791	1796	rVV5	115488	217182	1.60%	0.414%
60	13.475	1796	1800	1806	rVB	121794	191323	1.41%	0.365%
61	13.569	1806	1816	1819	rBV2	119294	272554	2.00%	0.520%
62	13.604	1820	1822	1831	rVB4	73553	108261	0.80%	0.206%
63	13.704	1833	1839	1842	rBV4	162274	271381	2.00%	0.517%
64	13.739	1842	1845	1849	rVB2	133669	188579	1.39%	0.359%
65	13.986	1883	1887	1892	rVB	851718	1095682	8.06%	2.089%
66	14.051	1892	1898	1906	rBV	1413801	2078554	15.28%	3.962%
67	14.163	1912	1917	1921	rVB2	139816	198187	1.46%	0.378%
68	14.210	1921	1925	1930	rVB	468937	626386	4.61%	1.194%
69	14.298	1935	1940	1943	rVV5	71396	139596	1.03%	0.266%
70	14.451	1960	1966	1971	rVV6	125844	247229	1.82%	0.471%
71	14.545	1978	1982	1989	rVB3	66705	112315	0.83%	0.214%
72	14.610	1989	1993	1997	rBV2	114143	159661	1.17%	0.304%
73	14.651	1997	2000	2002	rVV2	98378	121667	0.89%	0.232%
74	14.692	2002	2007	2012	rVV	717932	1048147	7.71%	1.998%
75	14.739	2012	2015	2019	rVV4	69819	115064	0.85%	0.219%
76	14.780	2019	2022	2026	rVB	143131	165315	1.22%	0.315%

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 Title : SVOA CALIBRATION

77	14.833	2026	2031	2037	rBV	487439	696452	5.12%	1.328%
78	14.951	2043	2051	2055	rBV	334388	498290	3.66%	0.950%
79	15.057	2065	2069	2074	rVB	146975	188384	1.39%	0.359%
80	15.174	2084	2089	2096	rBV	106701	200567	1.47%	0.382%
81	15.357	2114	2120	2125	rVB	98948	183611	1.35%	0.350%
82	15.445	2131	2135	2142	rVB7	58186	111793	0.82%	0.213%
83	15.716	2173	2181	2188	rBV2	265267	424355	3.12%	0.809%
84	15.816	2193	2198	2201	rBV	334045	478067	3.52%	0.911%
85	16.145	2251	2254	2264	rVV7	77400	163964	1.21%	0.313%
86	16.474	2303	2310	2321	rBV	9918956	13599802	100.00%	25.924%
87	16.616	2329	2334	2338	rBV	258099	342671	2.52%	0.653%
88	16.663	2339	2342	2351	rVB2	127903	224863	1.65%	0.429%
89	16.816	2361	2368	2378	rBV	1774643	2582485	18.99%	4.923%
90	16.910	2379	2384	2391	rVV	161870	302169	2.22%	0.576%
91	16.974	2391	2395	2400	rVB3	94259	126352	0.93%	0.241%
92	17.351	2452	2459	2464	rBV	241214	331661	2.44%	0.632%
93	17.527	2485	2489	2497	rVB2	111568	165535	1.22%	0.316%
94	18.045	2572	2577	2582	rBV	167917	211705	1.56%	0.404%
95	18.692	2682	2687	2689	rBV	144552	195362	1.44%	0.372%
96	19.221	2773	2777	2783	rVB2	174453	228415	1.68%	0.435%
97	19.821	2875	2879	2884	rVB3	107238	143392	1.05%	0.273%
98	20.792	3041	3044	3047	rBV	147060	154882	1.14%	0.295%
99	21.045	3082	3087	3096	rBV	2047115	2734150	20.10%	5.212%
100	23.756	3541	3548	3556	rVB	1209747	2649700	19.48%	5.051%

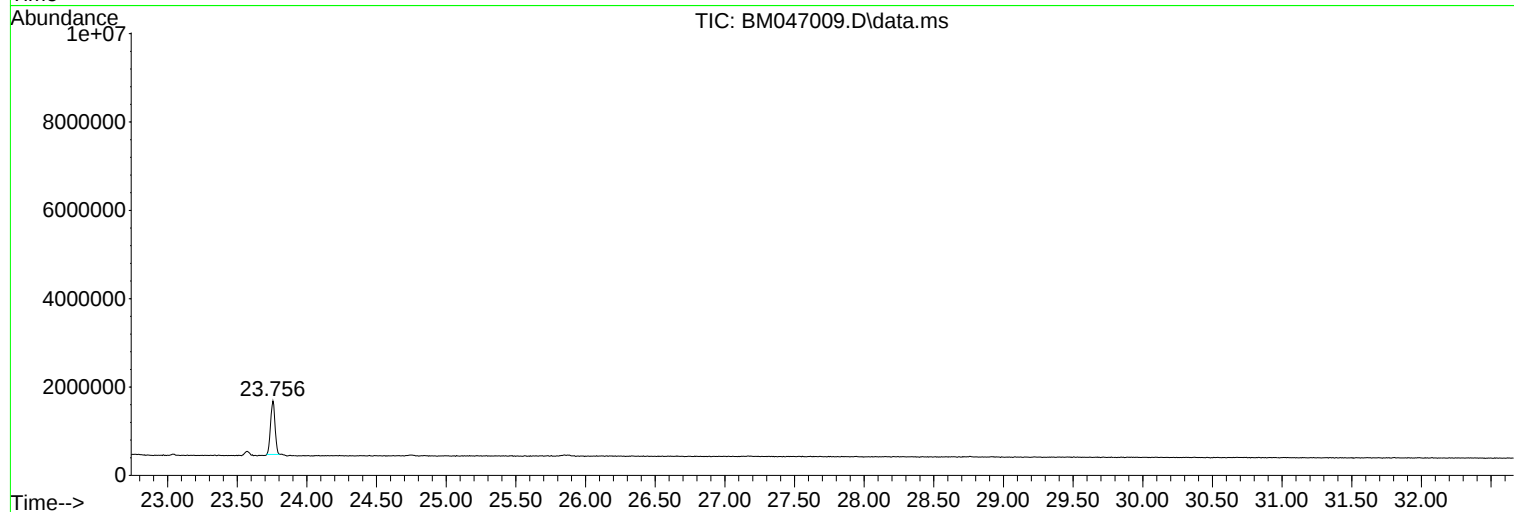
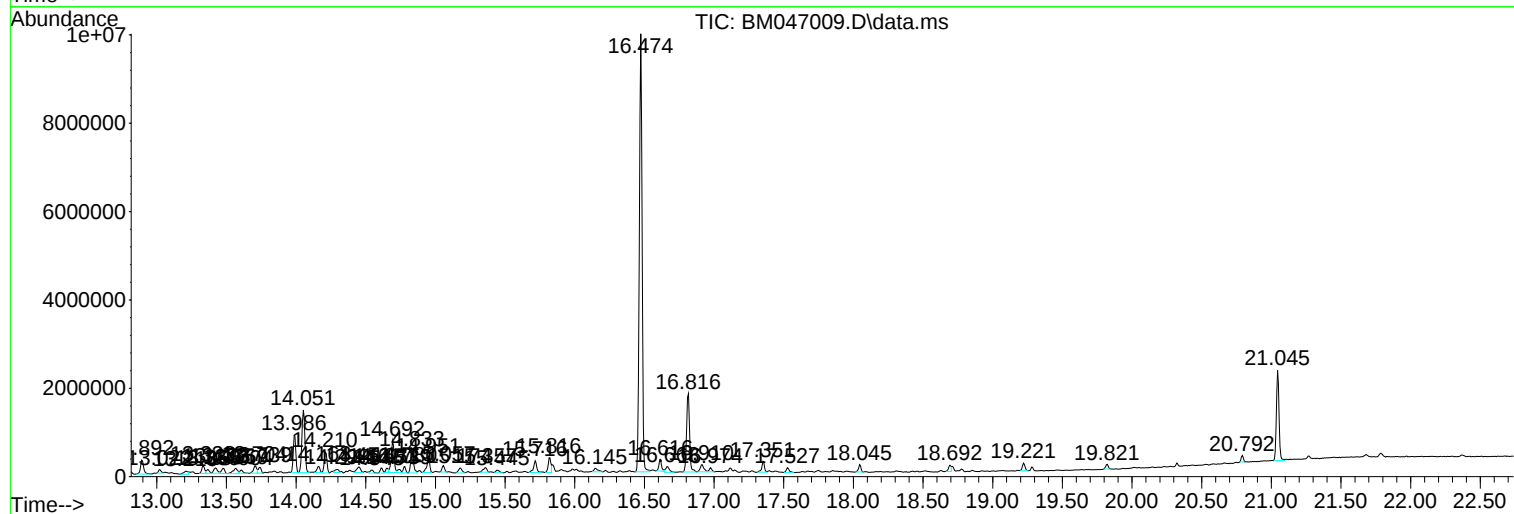
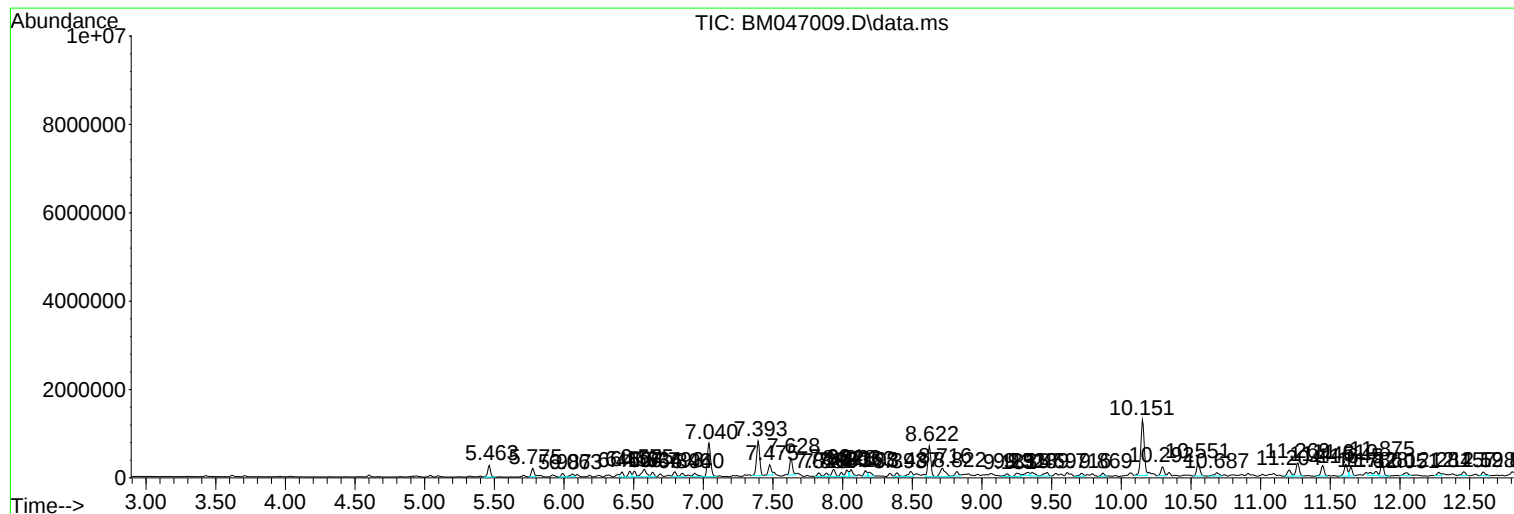
Sum of corrected areas: 52459934

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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



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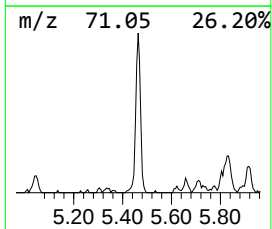
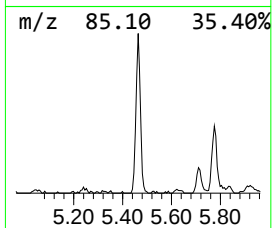
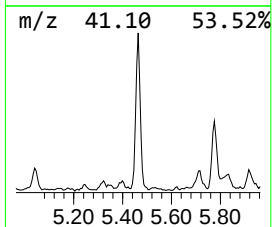
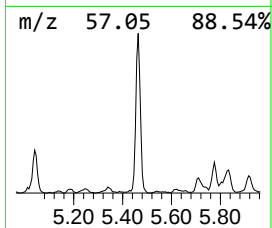
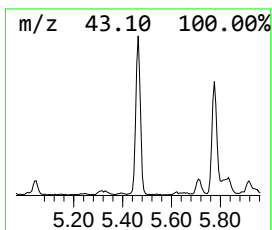
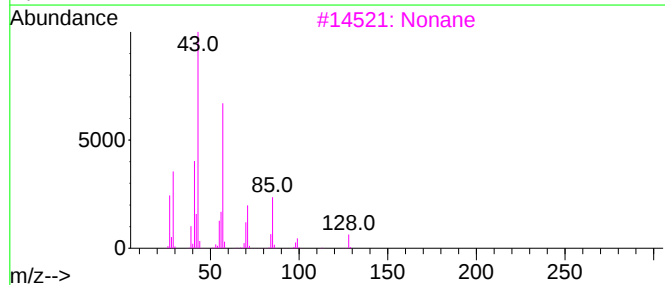
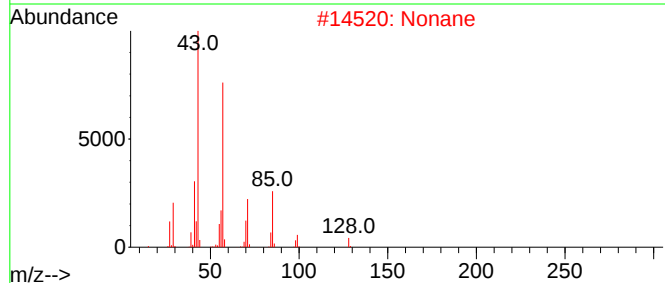
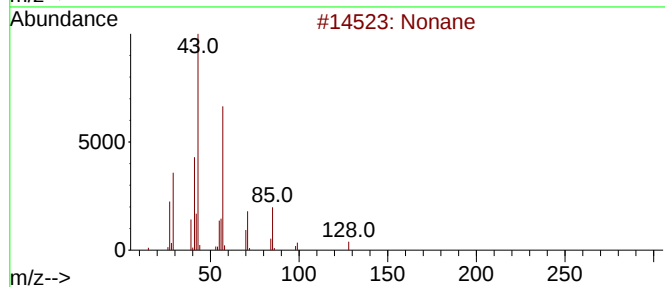
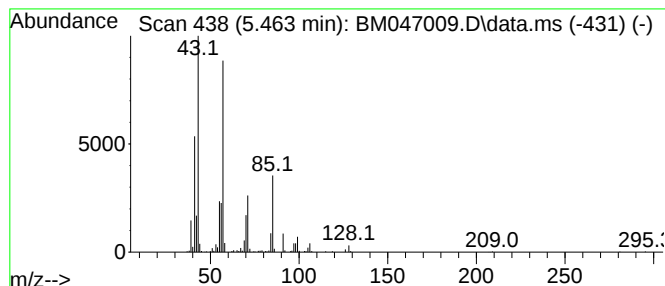
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.463	6.26 ng/ul	377853	1,4-Dichlorobenzene-d4	7.398

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	93
2	Nonane		128	C9H20	000111-84-2	90
3	Nonane		128	C9H20	000111-84-2	78
4	1-Octanol, 2-butyl-		186	C12H26O	003913-02-8	64
5	Decane		142	C10H22	000124-18-5	64



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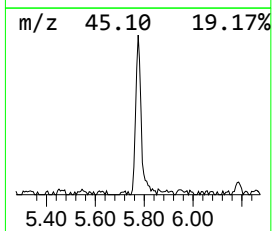
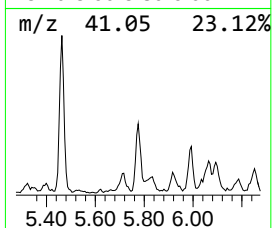
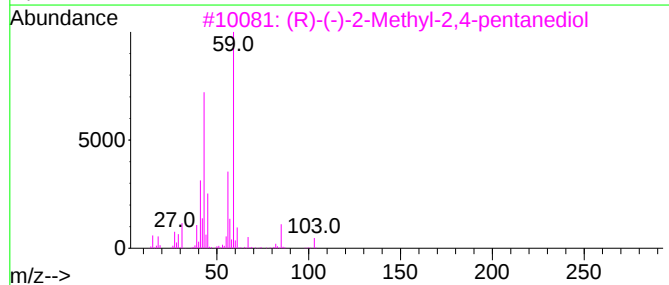
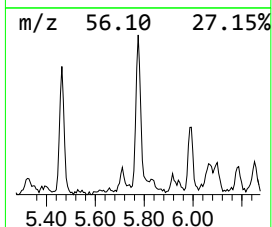
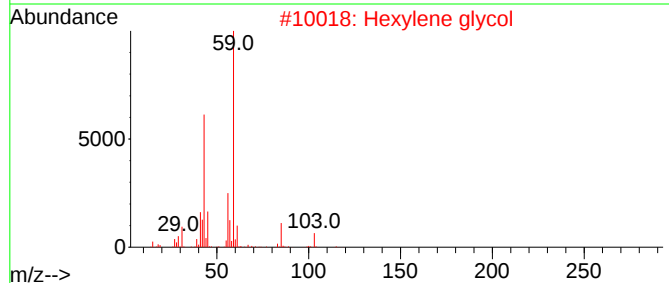
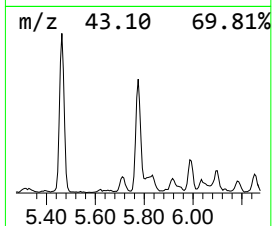
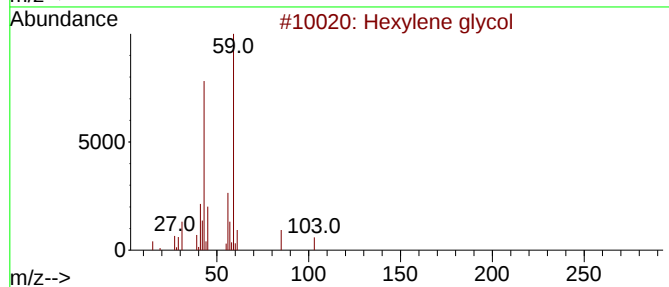
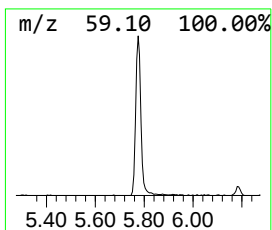
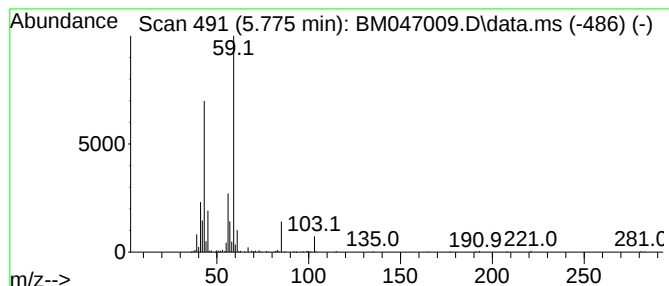
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.775	4.39 ng/ul	264976	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	90
3			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	90
4			Hexylene glycol	118	C6H14O2	000107-41-5	83
5			Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	78



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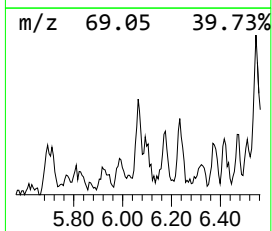
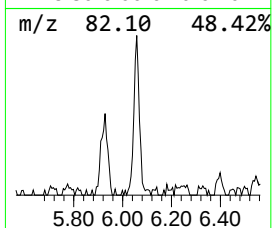
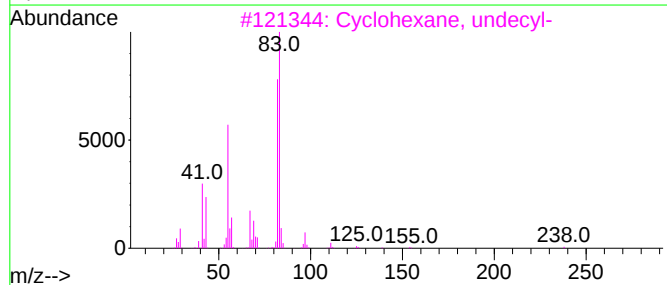
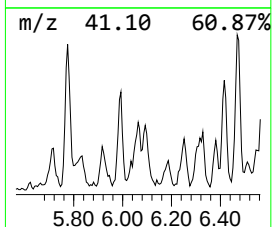
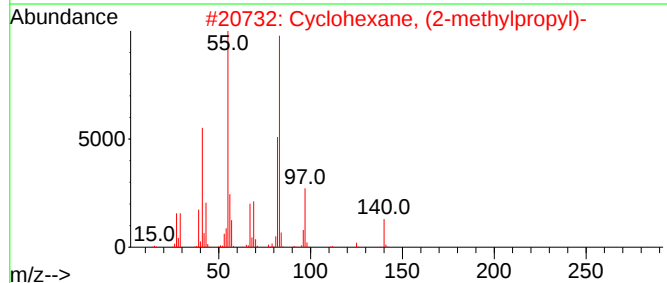
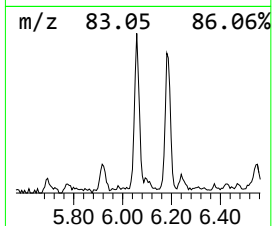
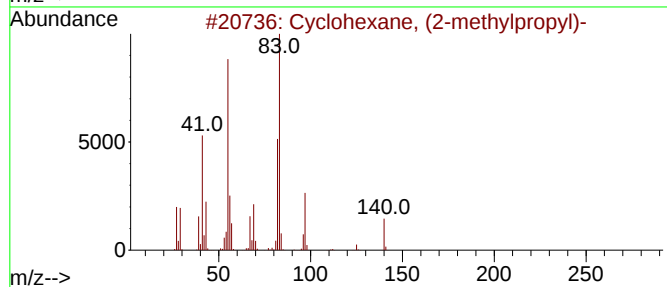
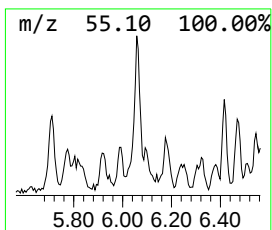
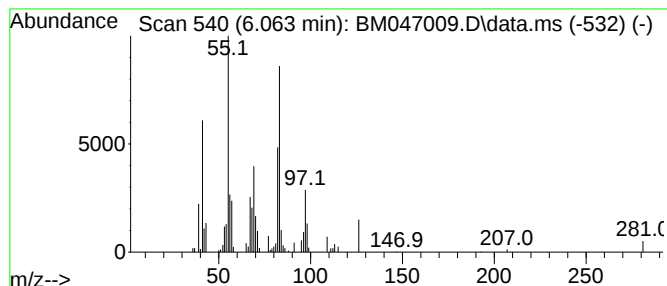
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic6.063 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.063	2.21 ng/ul	133588	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	59
2			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	59
3			Cyclohexane, undecyl-	238	C17H34	054105-66-7	53
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	52
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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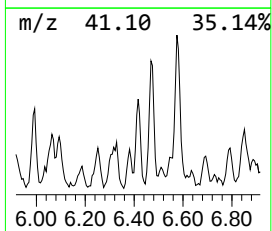
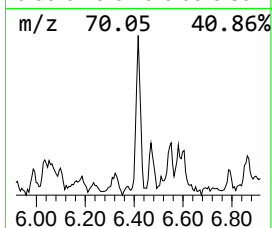
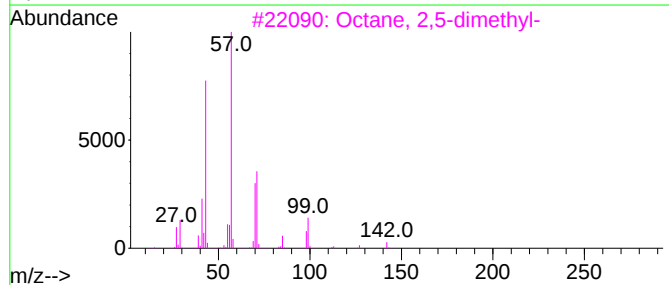
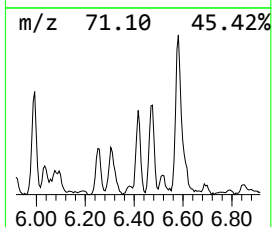
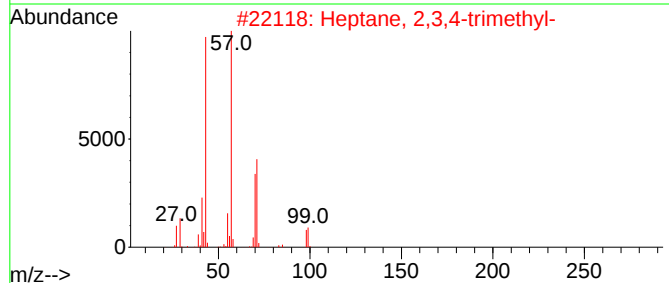
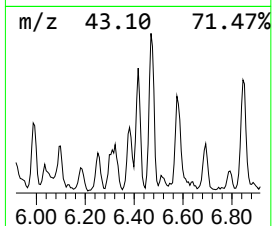
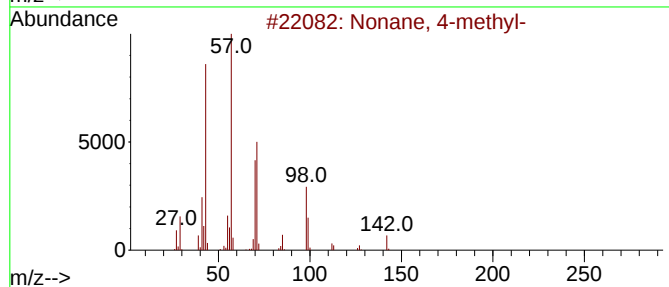
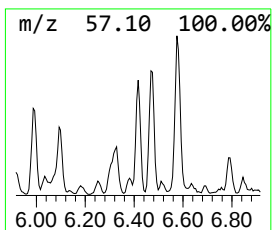
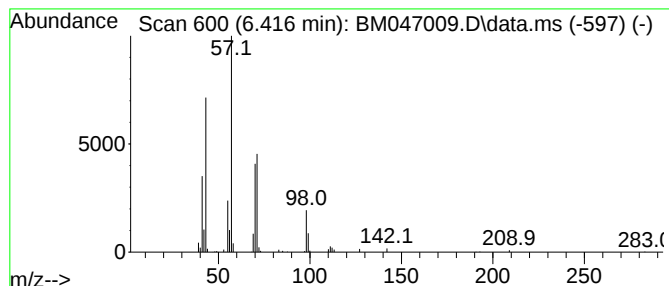
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.416	2.45 ng/ul	147650	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	83
2			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	72
3			Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	72
4			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	64
5			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
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Sample : P3171-05DL 30X
Misc :
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ClientSampleId :
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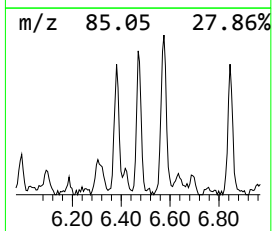
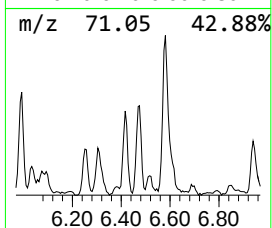
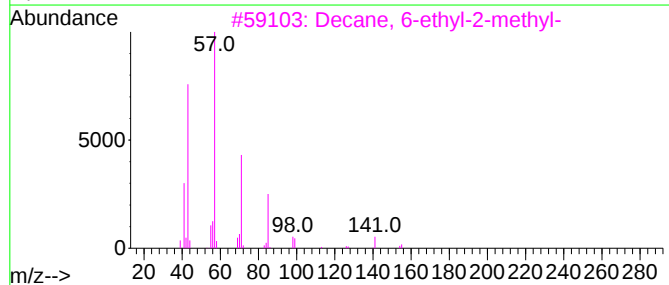
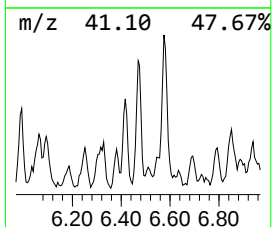
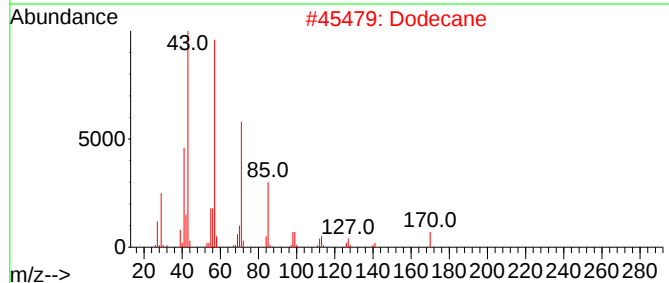
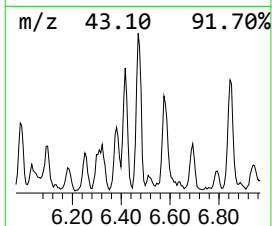
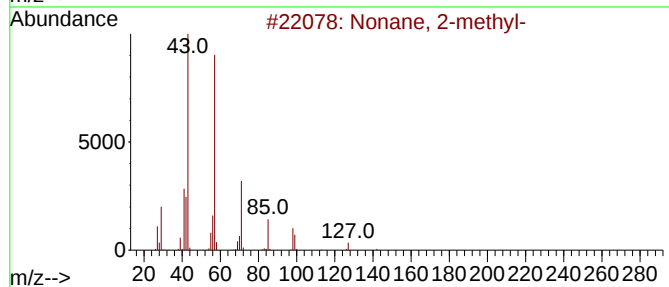
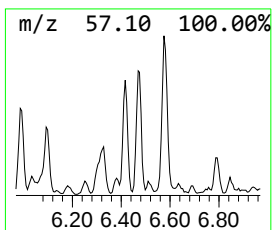
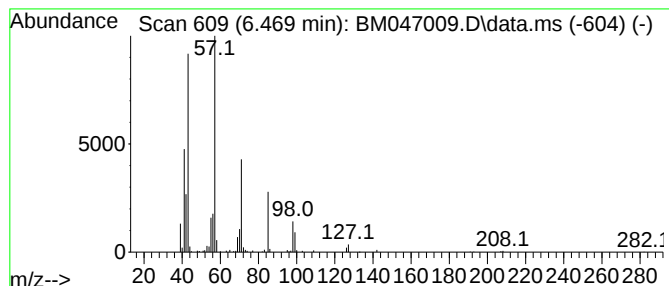
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.469	3.10 ng/ul	187087	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	90
2			Dodecane	170	C12H26	000112-40-3	72
3			Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64
4			Hexadecane	226	C16H34	000544-76-3	64
5			Nonane, 4-methyl-	142	C10H22	017301-94-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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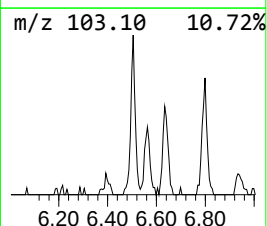
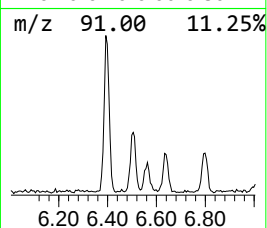
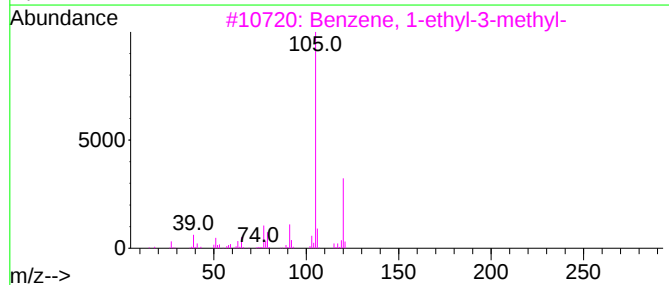
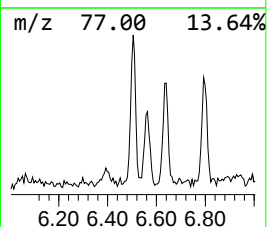
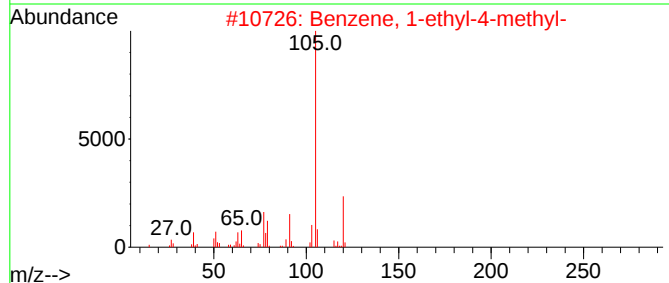
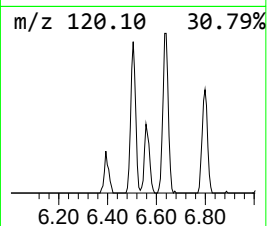
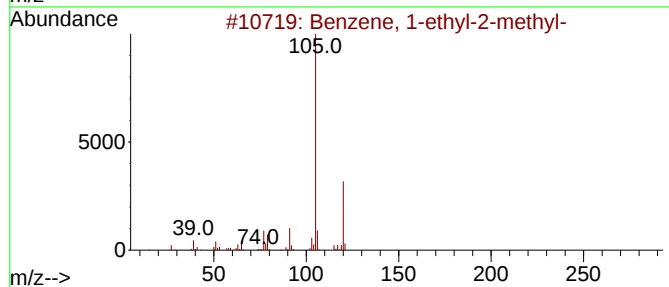
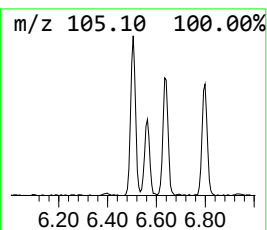
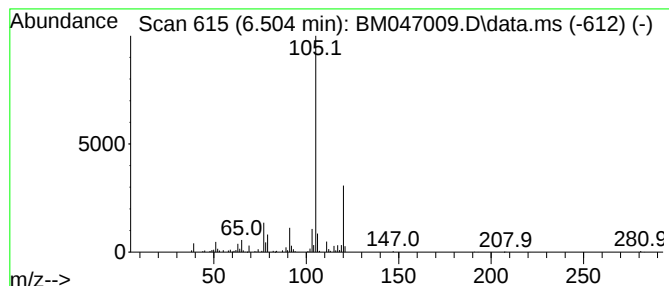
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.504	3.06 ng/ul	184651	1,4-Dichlorobenzene-d4	7.398

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-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
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ClientSampleId :
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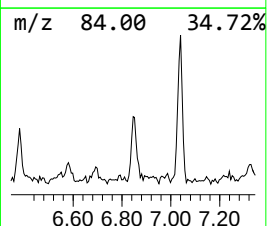
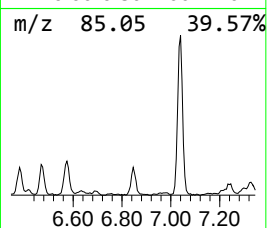
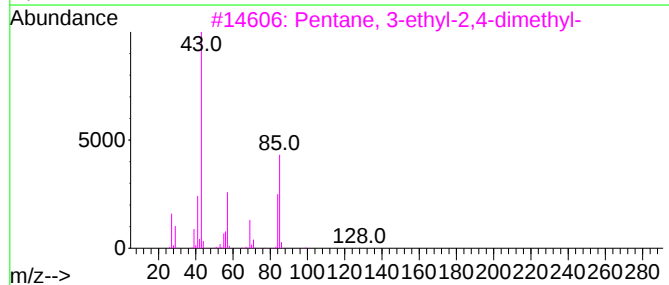
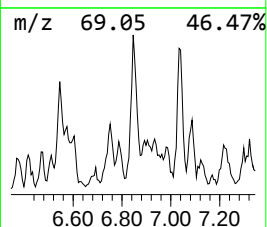
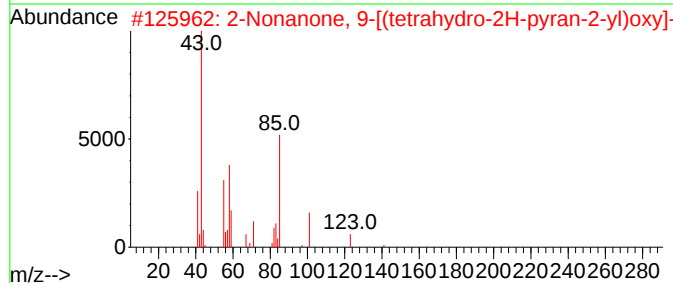
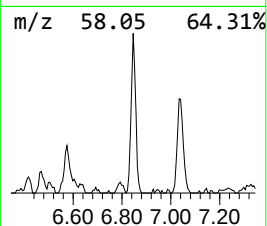
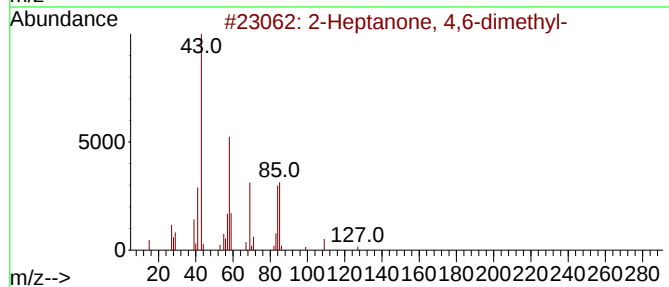
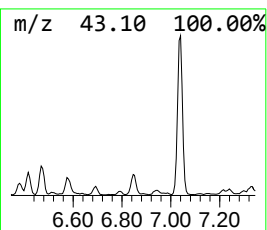
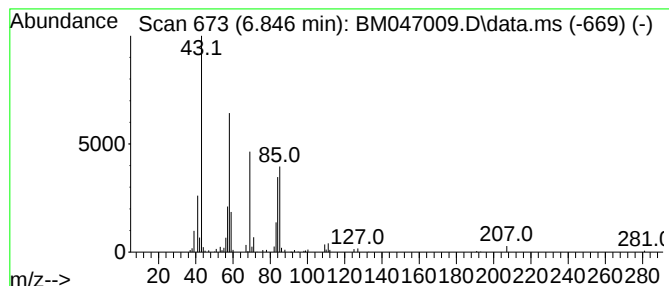
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-Heptanone, 4,6-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.846	2.36 ng/ul	142251	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	83
2			2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	50
3			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	43
4			Hexane, 3-ethyl-	114	C8H18	000619-99-8	38
5			2-Pentanone, 5,5'-oxybis-	186	C10H18O3	093677-66-8	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
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Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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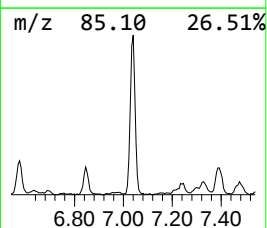
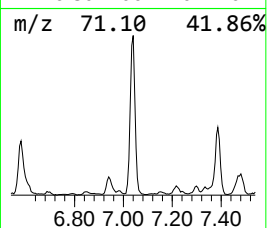
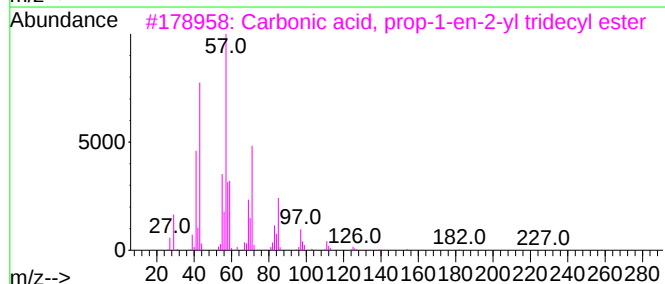
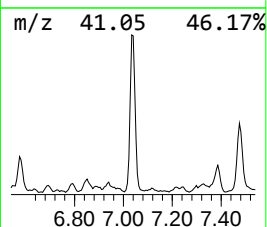
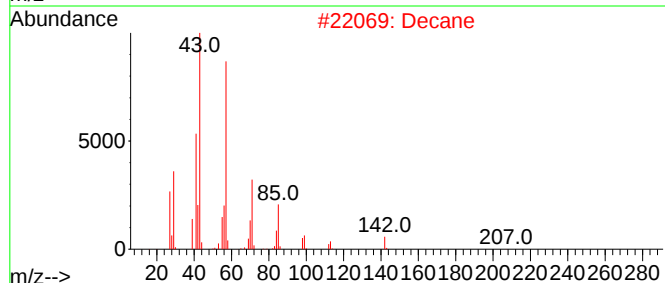
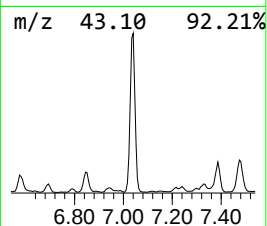
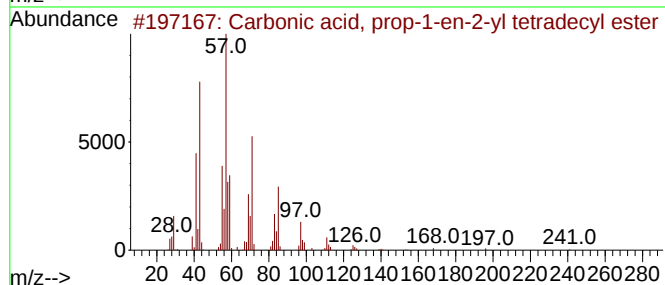
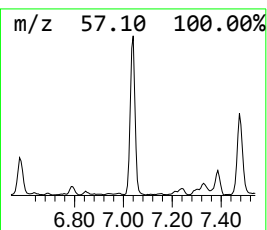
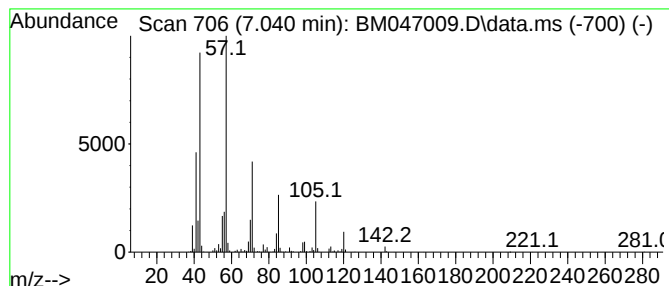
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Carbonic acid, prop-1-en-2-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.040	20.58 ng/ul	1242900	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	64
2			Decane	142	C10H22	000124-18-5	64
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	59
4			Decane	142	C10H22	000124-18-5	55
5			Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId :
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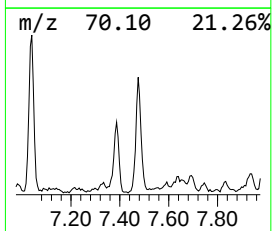
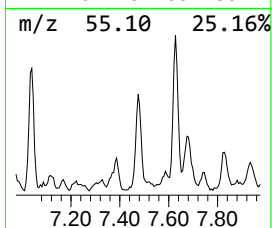
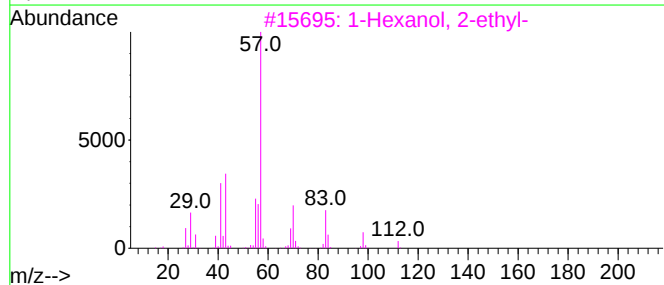
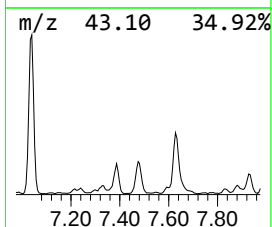
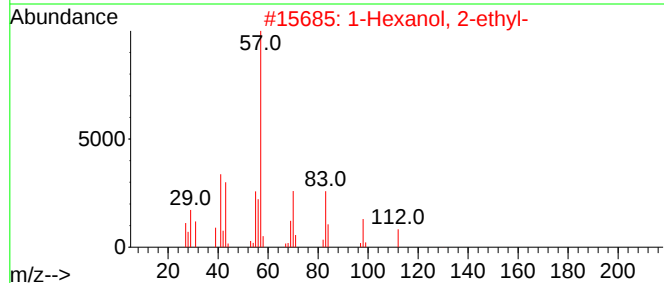
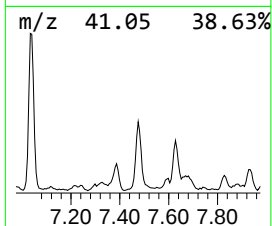
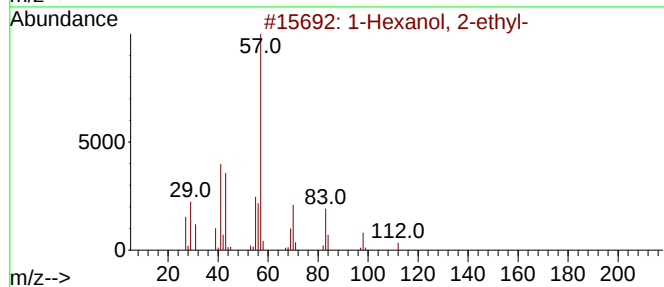
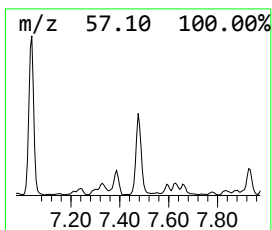
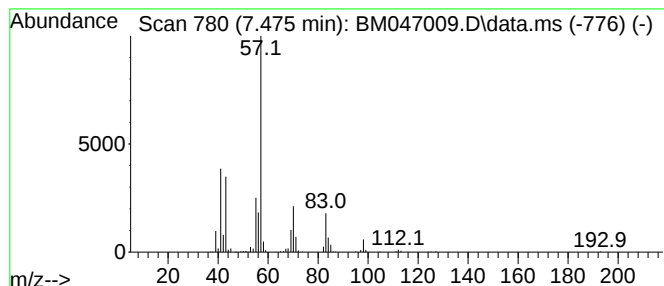
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Hexanol, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.475	5.85 ng/ul	353087	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	Heptane, 1,1'-oxybis-			214	C14H30O	000629-64-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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BNA_M
ClientSampleId :
DCVW9DL

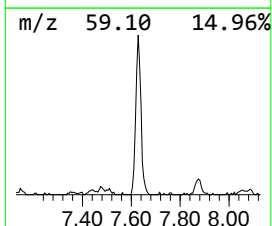
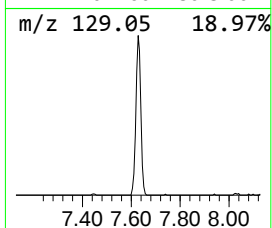
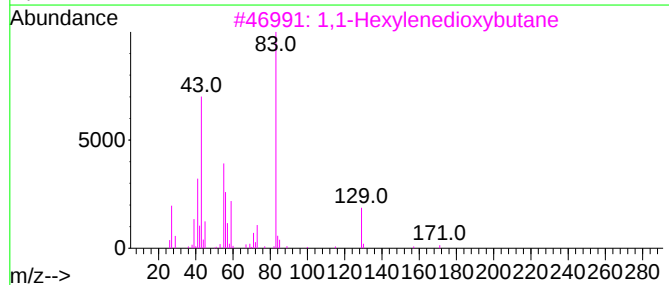
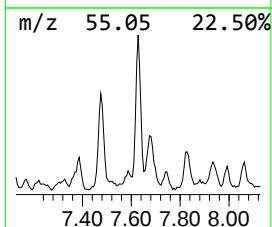
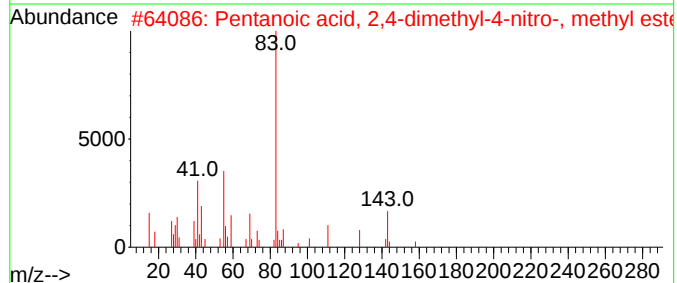
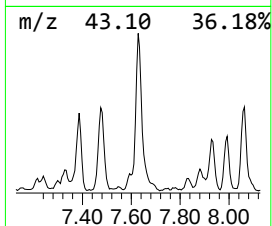
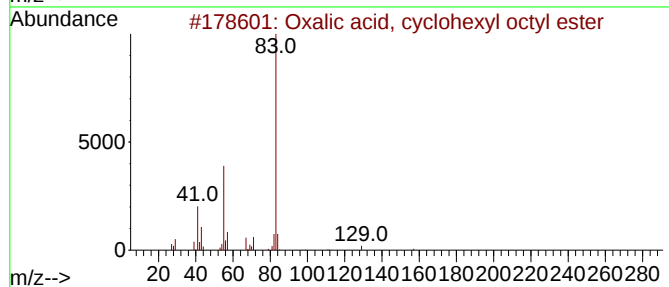
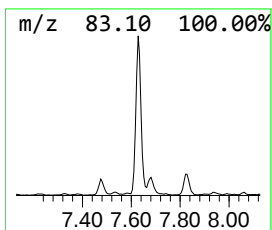
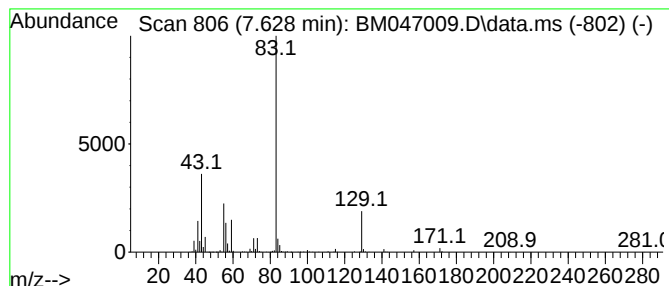
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Oxalic acid, cyclohexyl oct... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.628	9.28 ng/ul	560356	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	53
2			Pentanoic acid, 2,4-dimethyl-4-n...	189	C8H15NO4	005762-40-3	40
3			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	38
4			Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	36
5			Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

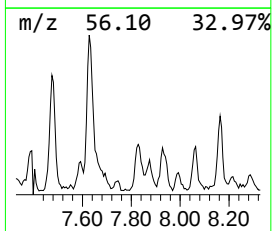
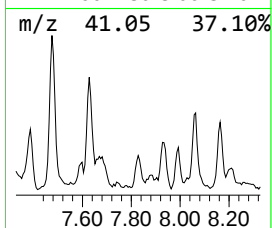
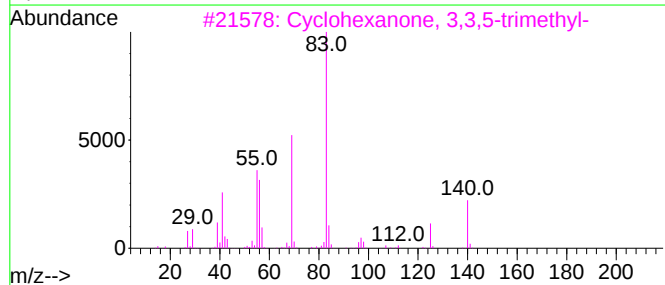
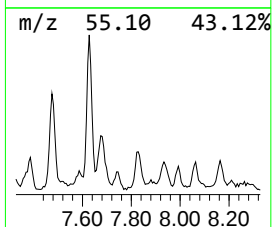
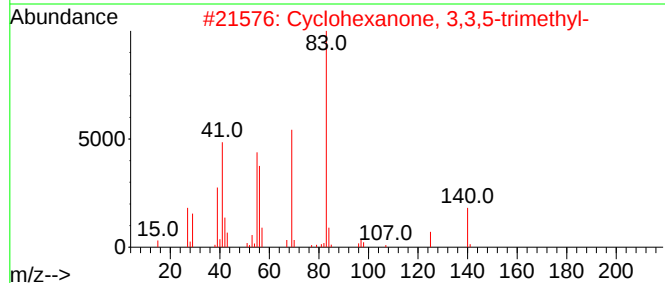
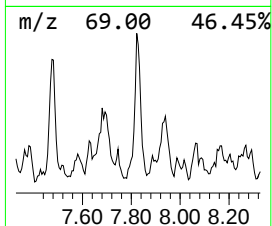
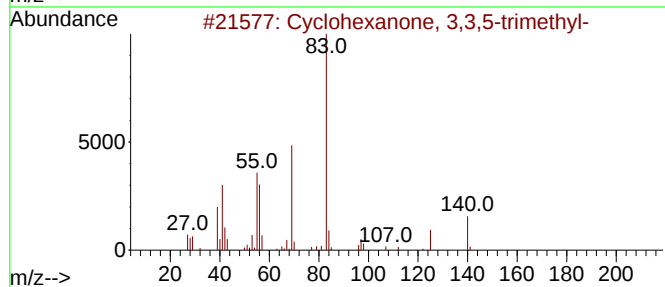
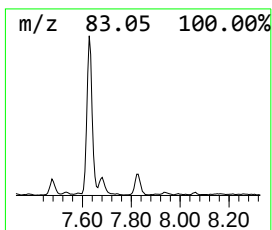
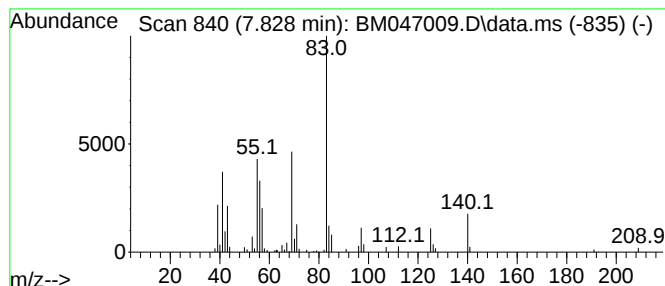
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Cyclohexanone, 3,3,5-trimet... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	2.15 ng/ul	130017	1,4-Dichlorobenzene-d4	7.398

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
2		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	64
5		1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

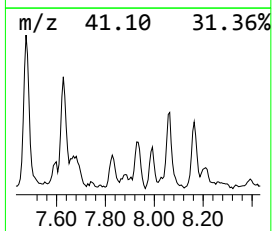
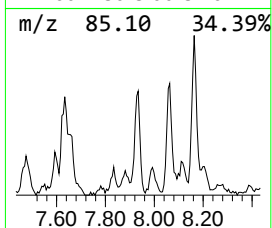
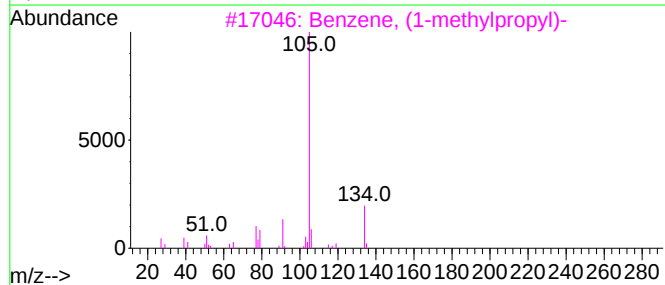
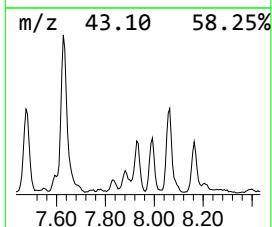
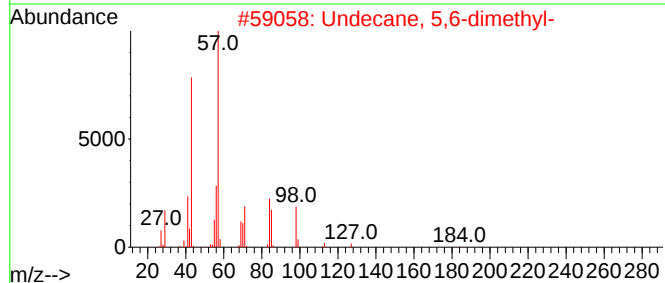
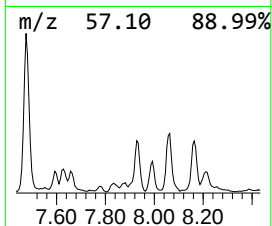
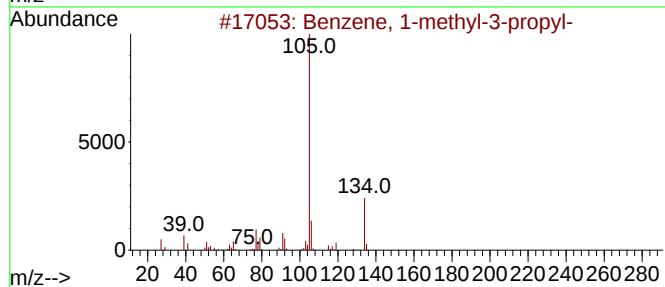
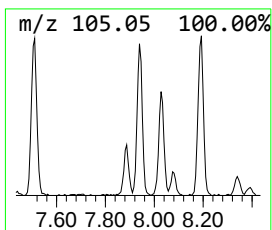
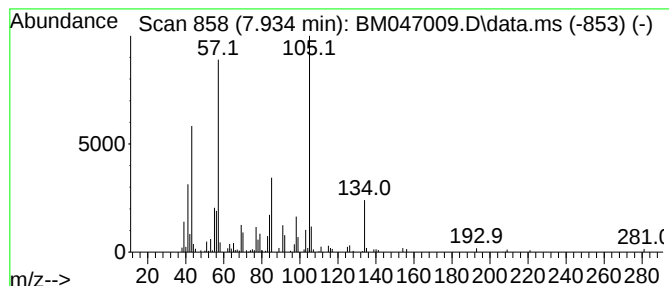
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	4.45 ng/ul	268914	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
2			Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	38
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	38
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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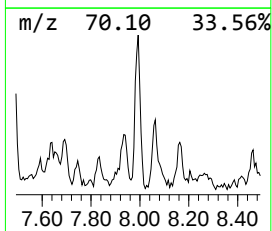
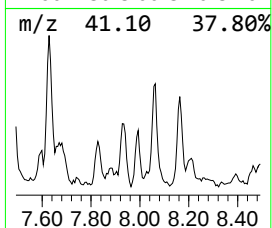
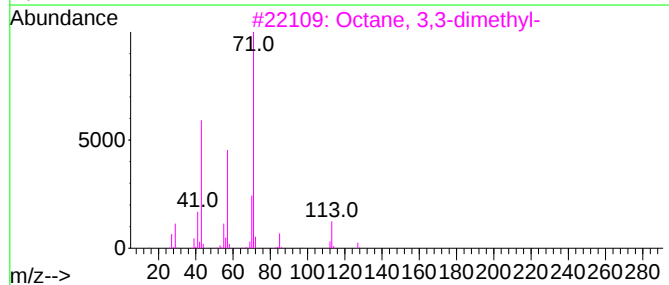
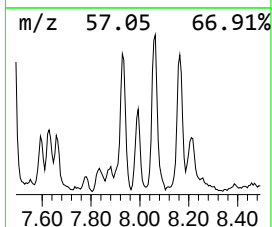
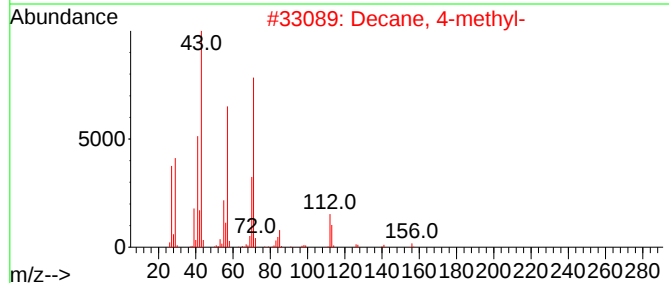
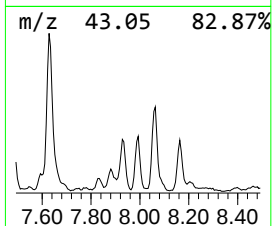
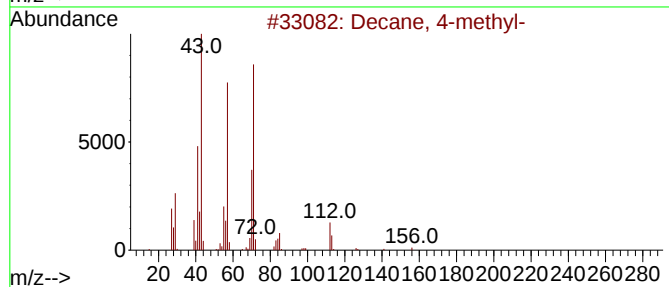
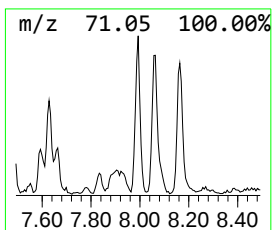
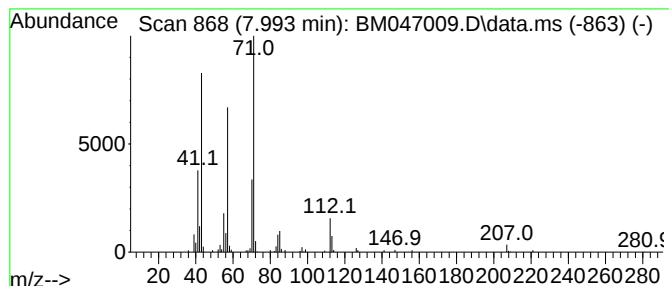
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	2.10 ng/ul	126638	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	91
2			Decane, 4-methyl-	156	C11H24	002847-72-5	83
3			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
4			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	78
5			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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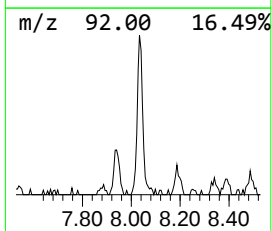
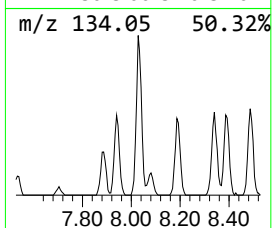
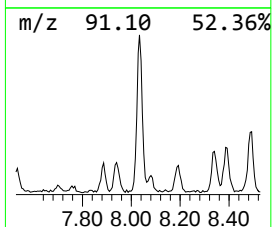
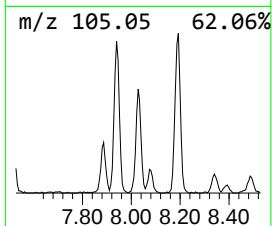
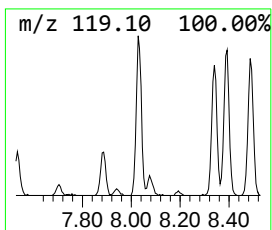
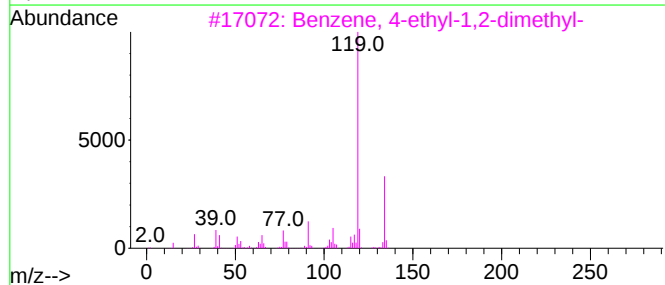
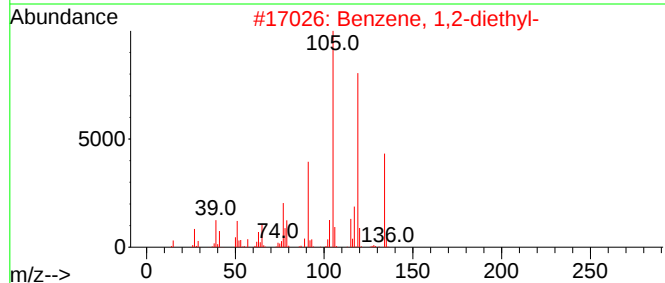
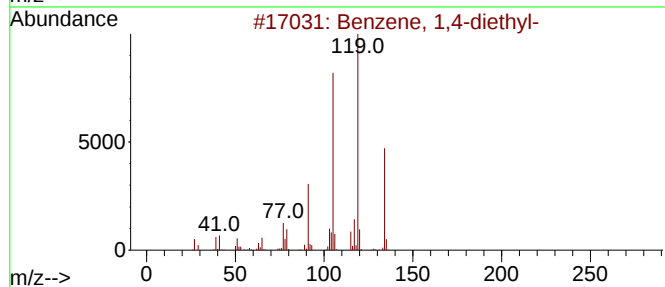
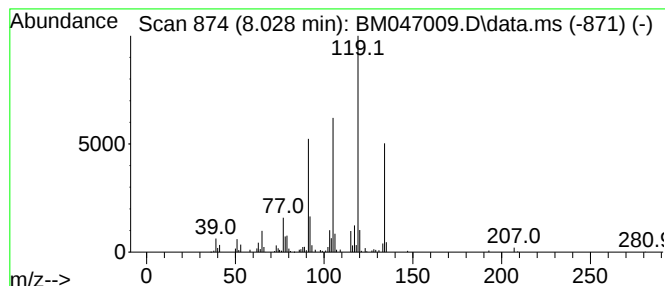
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1,4-diethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.028	3.54 ng/ul	213831	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

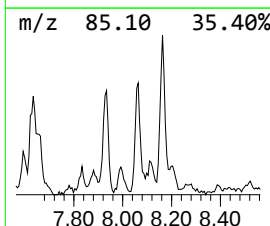
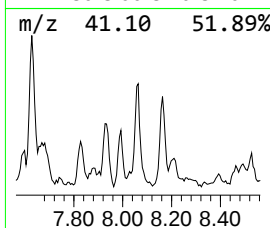
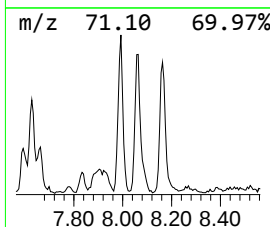
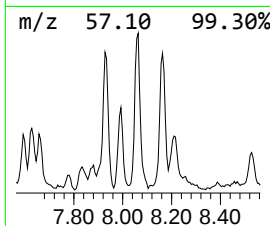
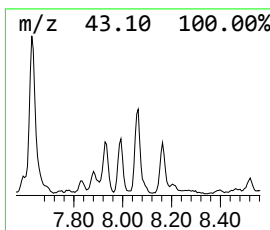
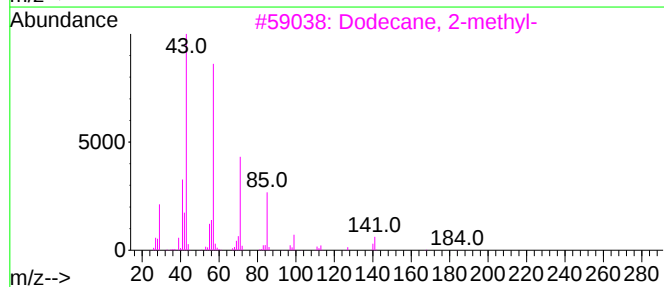
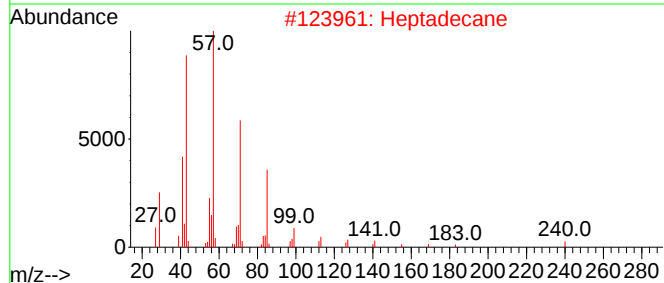
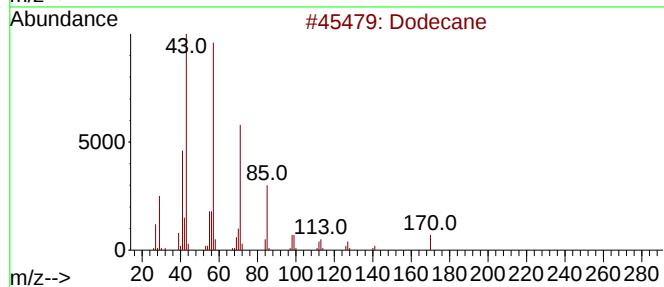
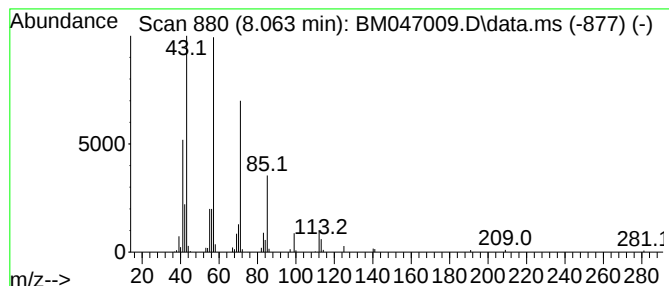
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.063	3.77 ng/ul	227787	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	83
2			Heptadecane	240	C17H36	000629-78-7	72
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
4			Tridecane	184	C13H28	000629-50-5	72
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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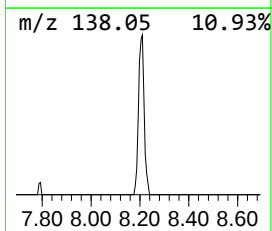
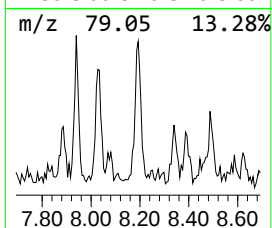
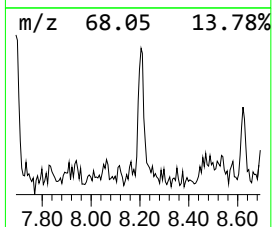
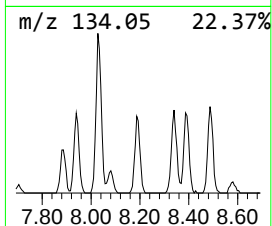
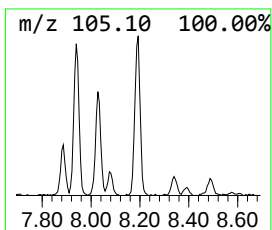
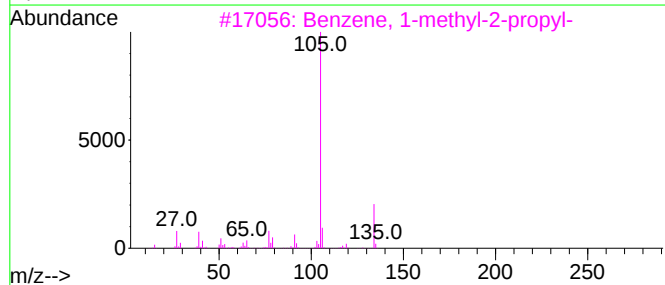
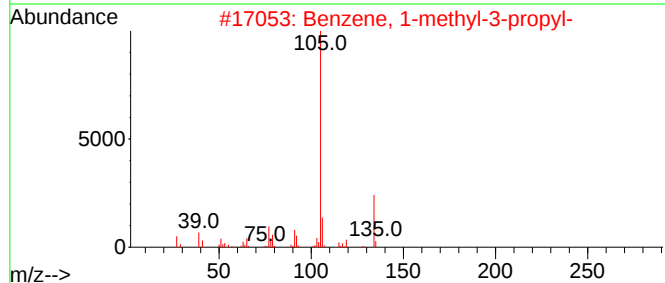
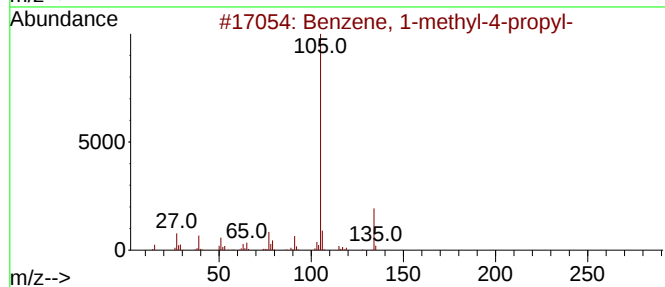
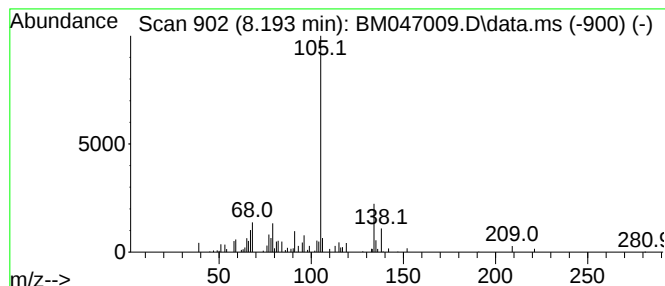
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzene, 1-methyl-4-propyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.193	2.43 ng/ul	147023	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	52
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	52
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	52
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	50
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	50



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Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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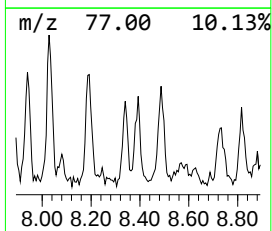
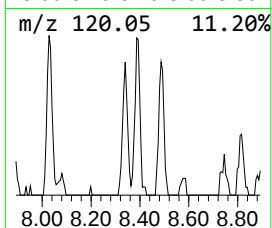
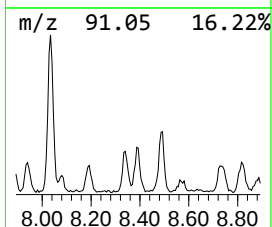
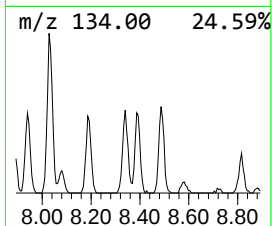
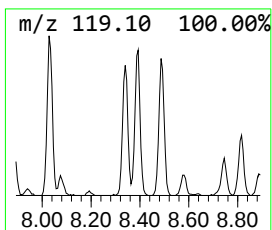
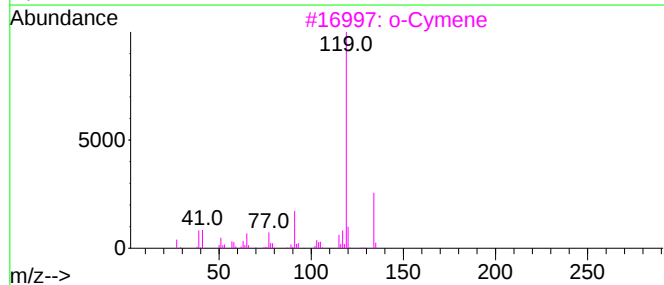
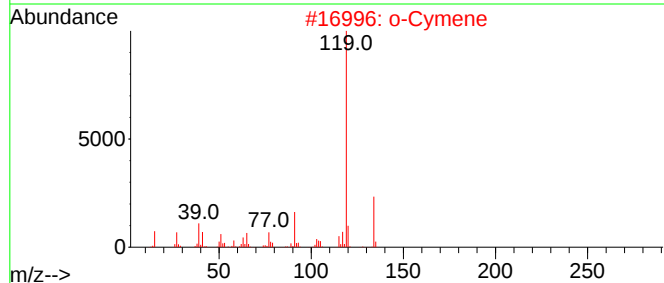
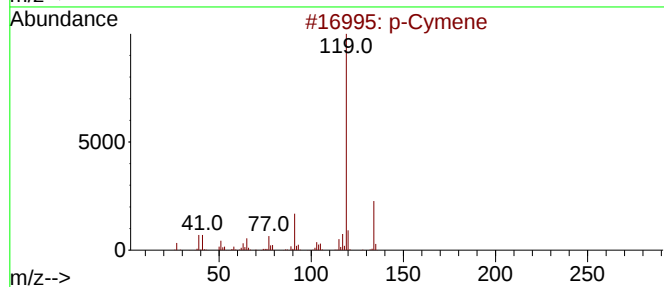
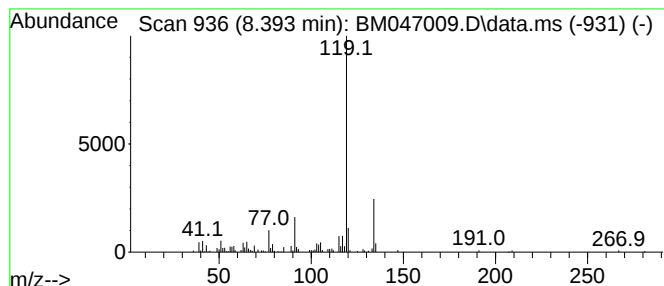
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 p-Cymene Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.393	2.08 ng/ul	125303	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			o-Cymene	134	C10H14	000527-84-4	97
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

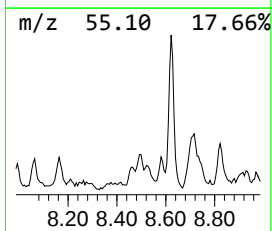
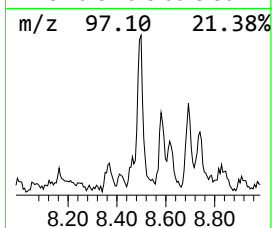
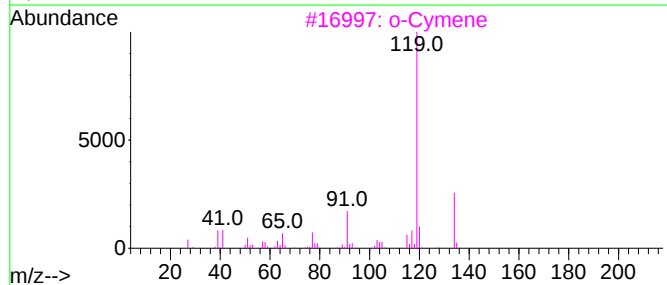
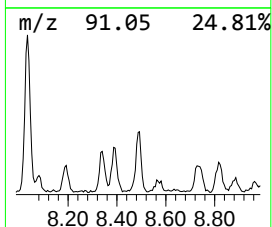
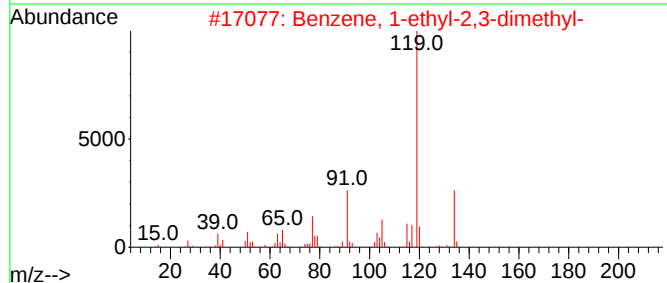
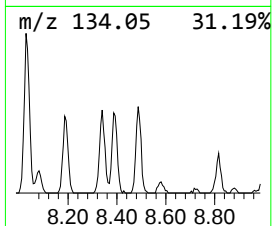
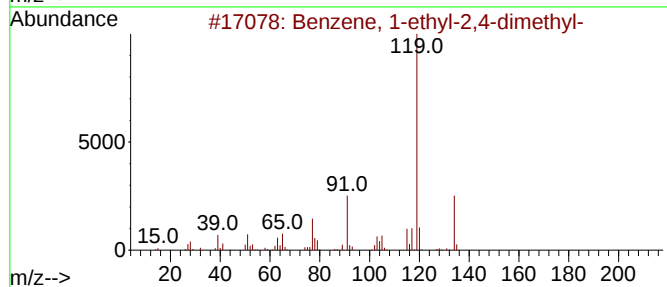
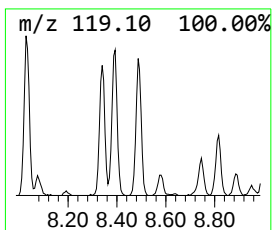
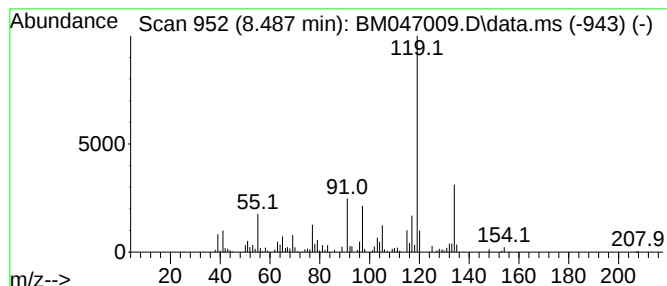
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.487	3.88 ng/ul	234229	1,4-Dichlorobenzene-d4	7.398

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	96
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3		o-Cymene	134	C10H14	000527-84-4	94
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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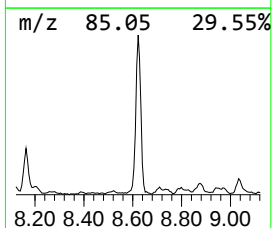
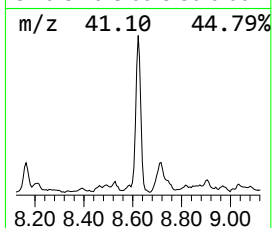
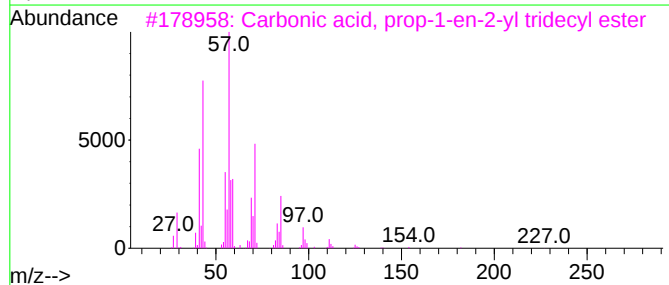
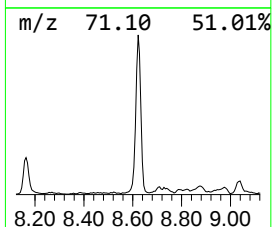
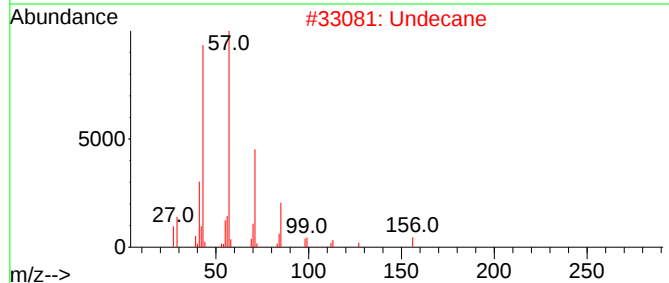
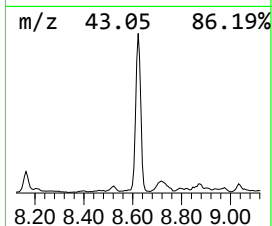
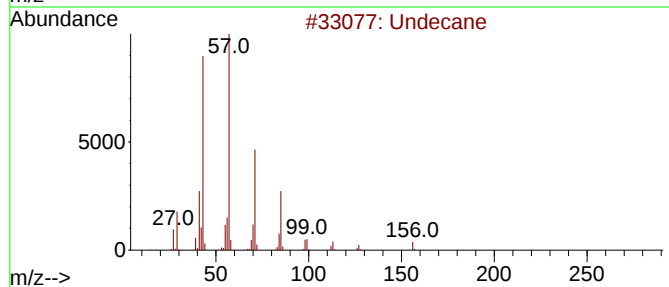
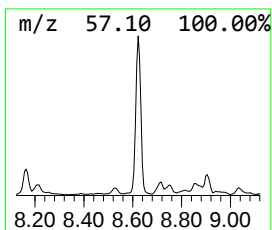
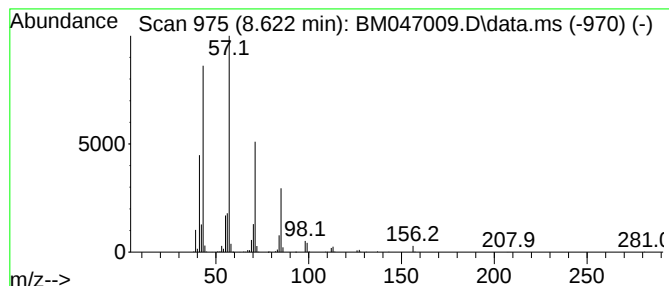
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.622	16.62 ng/ul	1003790	1,4-Dichlorobenzene-d4	7.398

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	Hexadecane			226	C16H34	000544-76-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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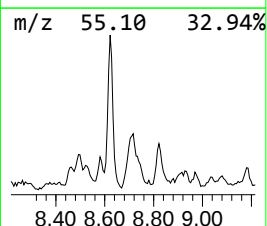
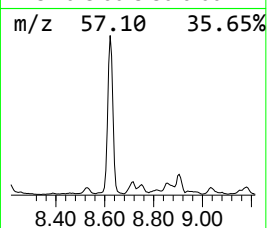
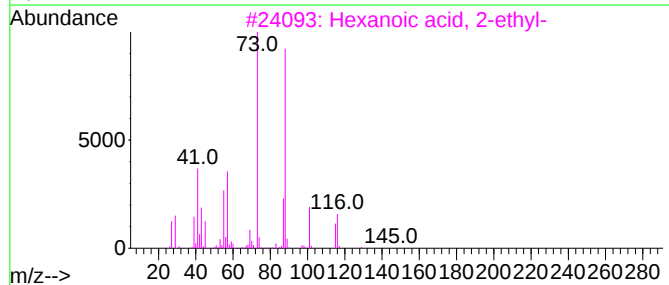
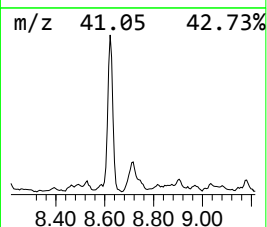
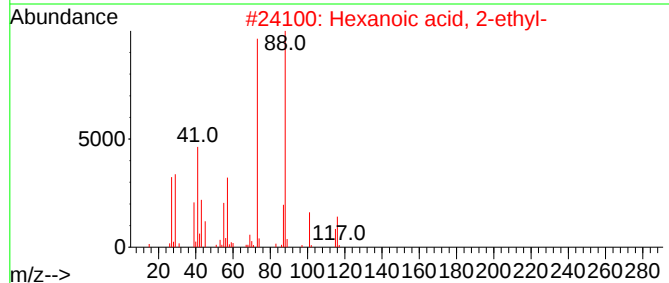
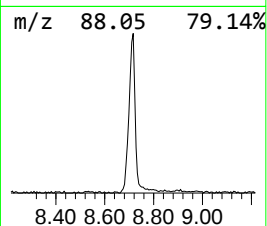
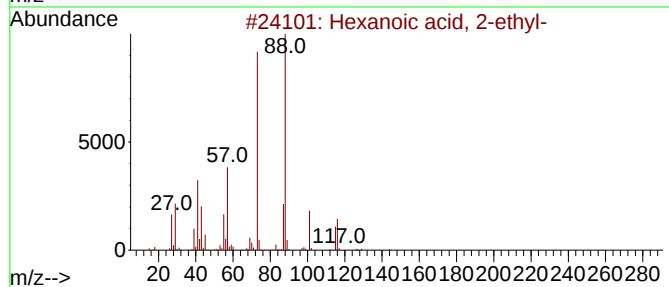
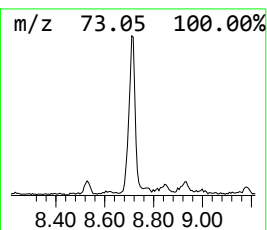
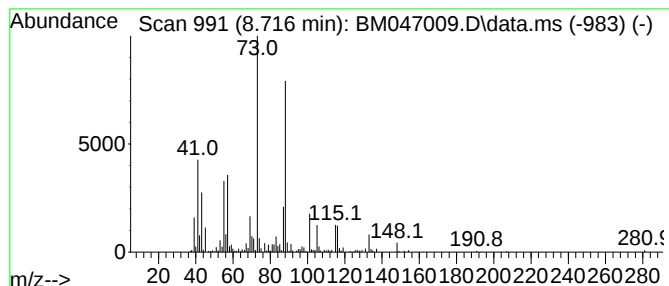
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Hexanoic acid, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.716	8.34 ng/ul	503381	1,4-Dichlorobenzene-d4	7.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
4			2-Ethyl-4-methylpentanoic acid	144	C8H16O2	000108-81-6	64
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
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ClientSampleId :
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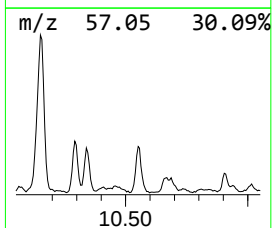
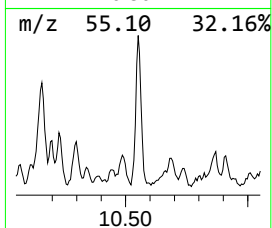
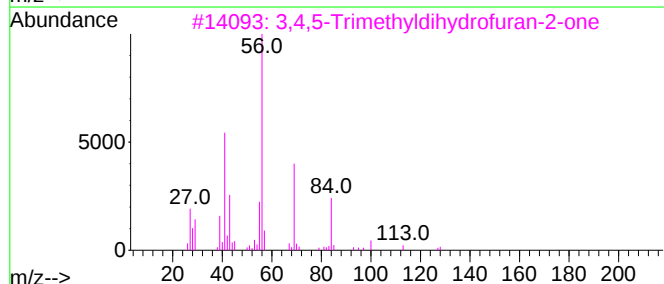
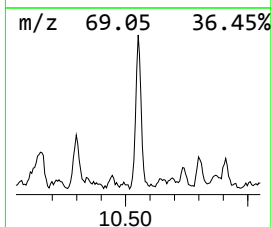
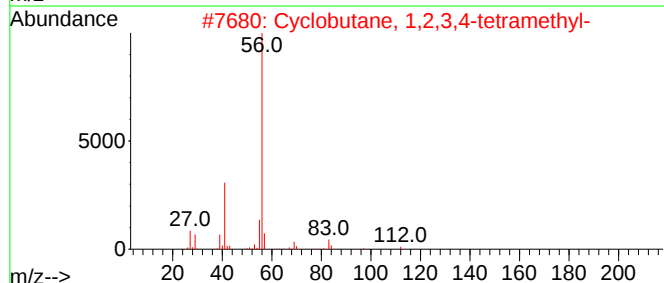
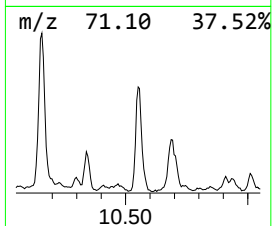
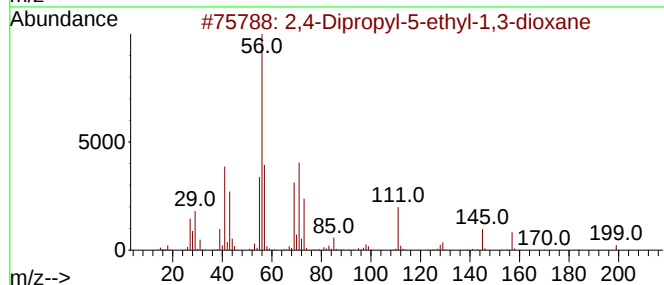
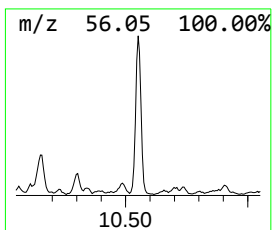
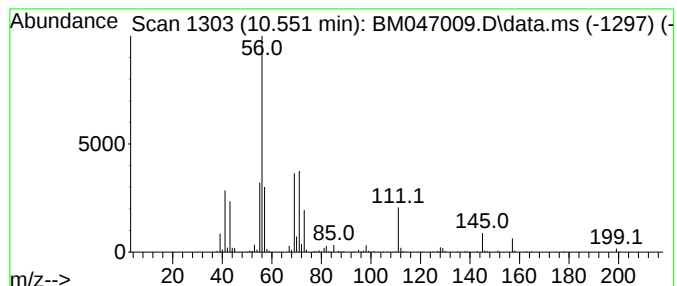
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.551	4.39 ng/ul	483162	Naphthalene-d8	10.151

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	47
2		Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	38
3		3,4,5-Trimethyldihydrofuran-2-one	128	C7H12O2	1000194-05-2	32
4		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	23
5		Cyclohexylamine	99	C6H13N	000108-91-8	23



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId :
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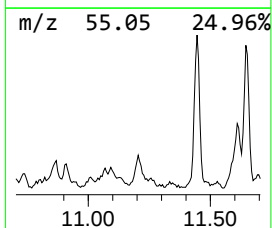
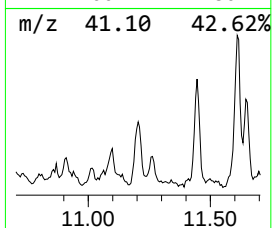
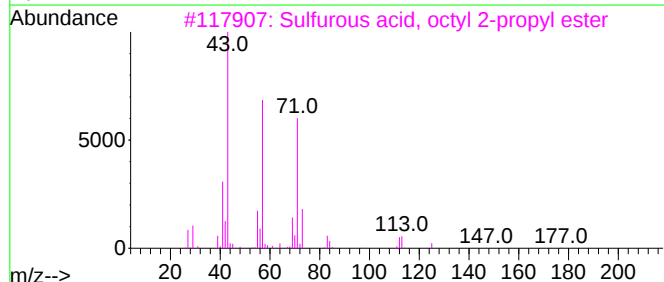
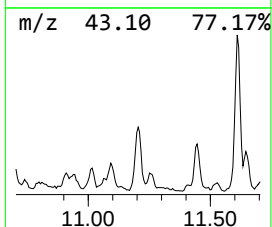
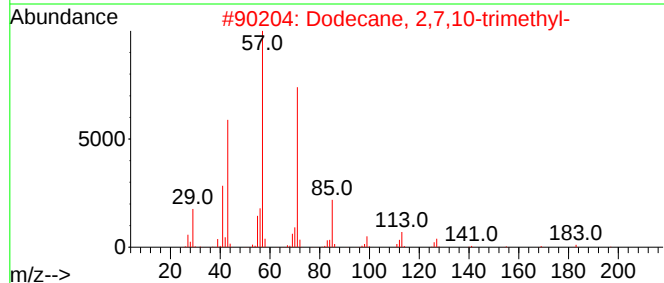
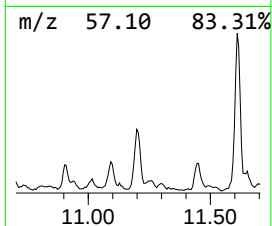
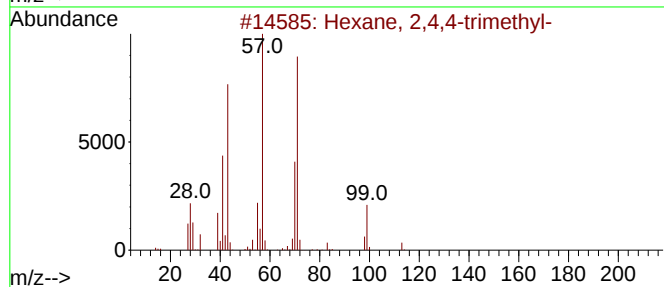
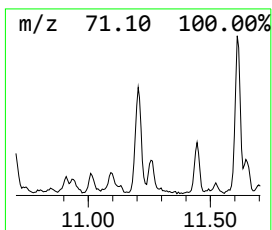
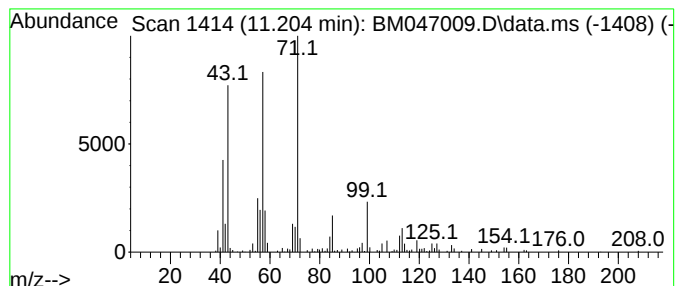
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.204	2.47 ng/ul	271738	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	53
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	53
3			Sulfurous acid, octyl 2-propyl e...	236	C11H24O3S	1000309-11-9	50
4			2,2,3-Triethyloxirane	128	C8H16O	053229-40-6	47
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
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Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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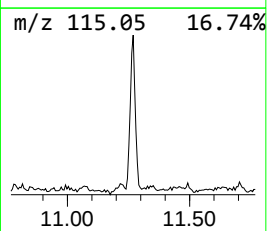
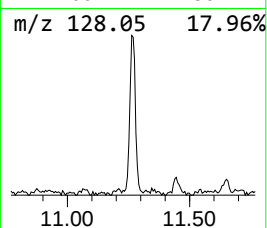
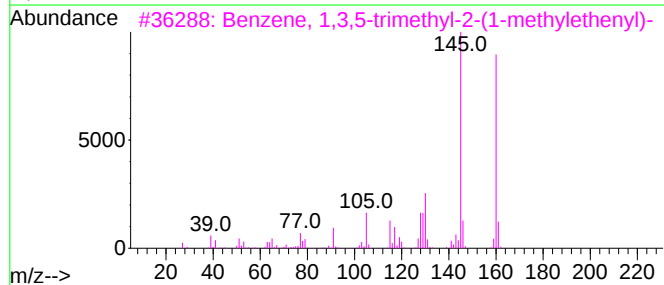
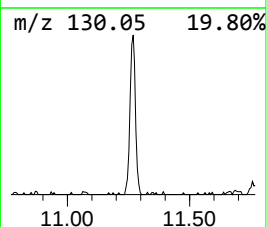
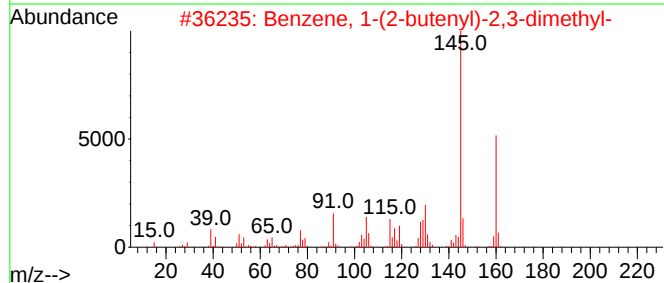
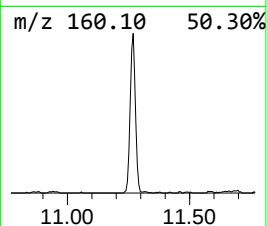
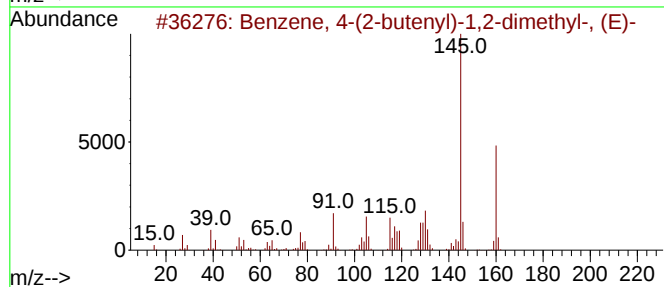
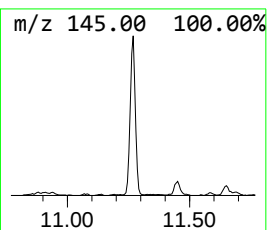
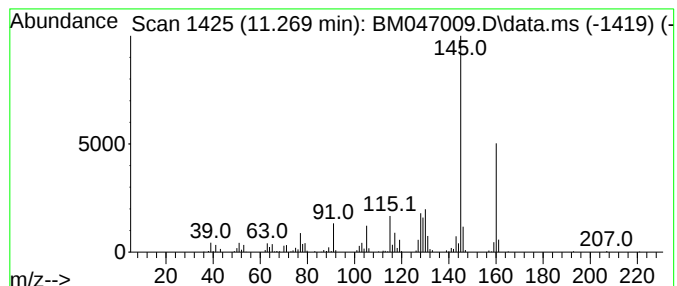
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.269	4.39 ng/ul	483654	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
2			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
3			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	93
4			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	91
5			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

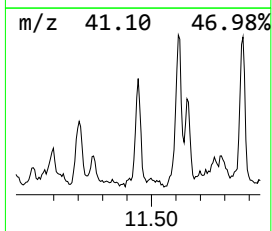
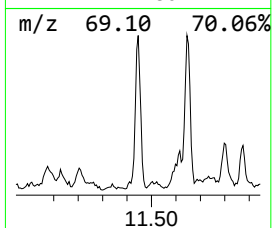
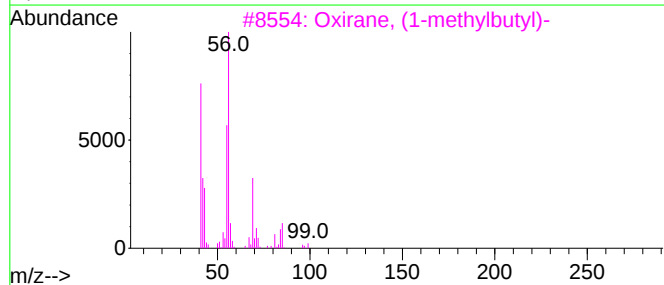
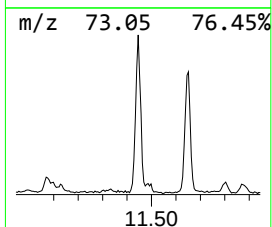
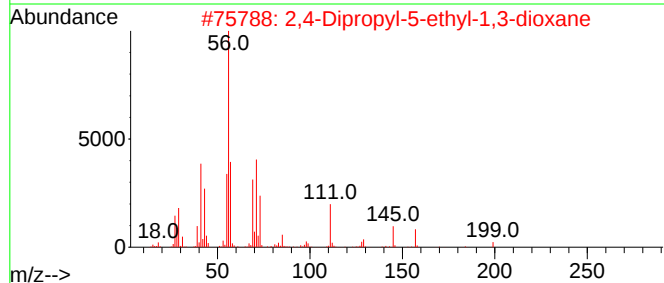
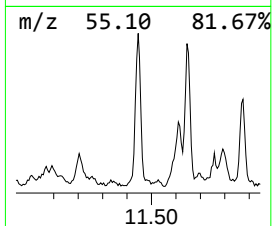
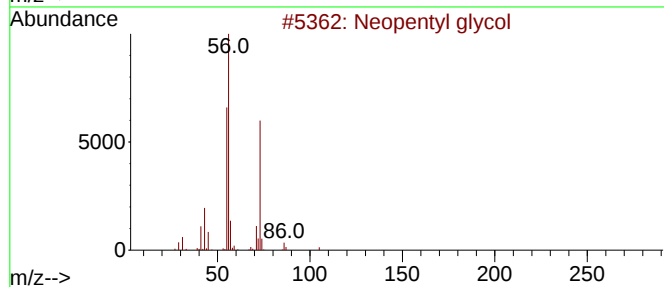
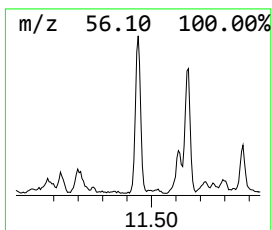
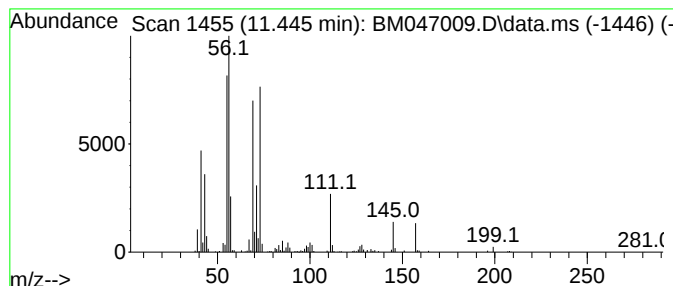
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 unknown-03 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.445	3.49 ng/ul	384599	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Neopentyl glycol	104	C5H12O2	000126-30-7	40
2			2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	37
3			Oxirane, (1-methylbutyl)-	114	C7H14O	053229-39-3	37
4			Pentanal, 3-methyl-	100	C6H12O	015877-57-3	35
5			Neopentyl glycol	104	C5H12O2	000126-30-7	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

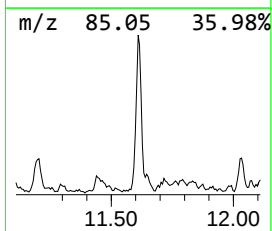
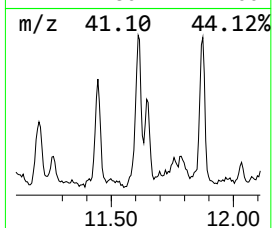
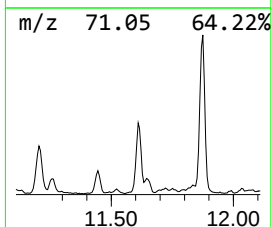
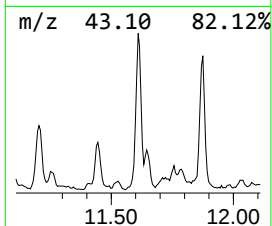
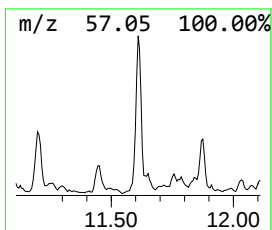
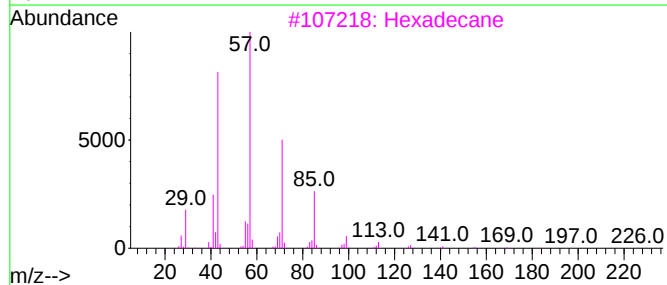
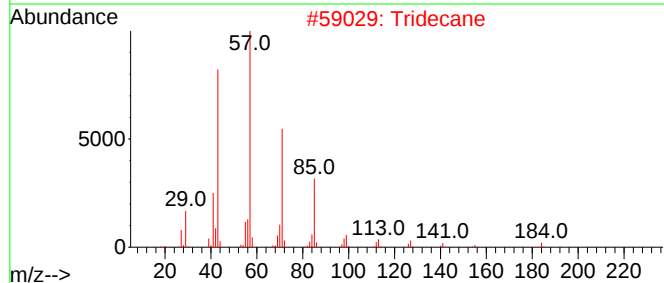
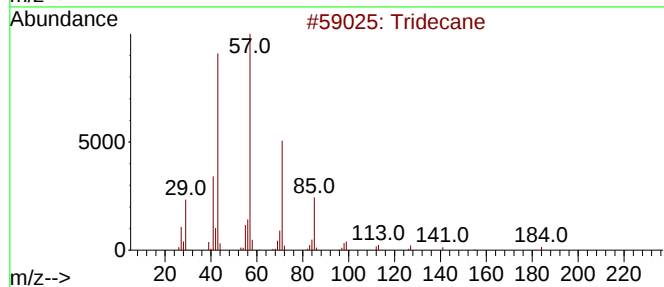
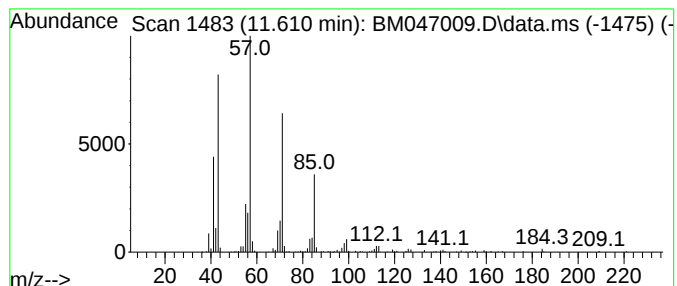
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	4.39 ng/ul	482791	Naphthalene-d8	10.151

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	93
2		Tridecane	184	C13H28	000629-50-5	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Tetradecane	198	C14H30	000629-59-4	86
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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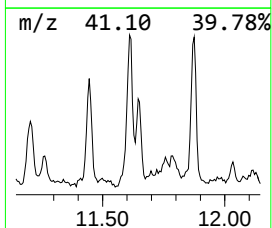
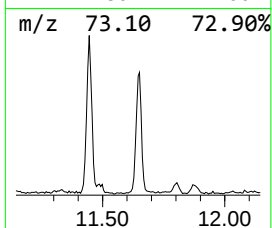
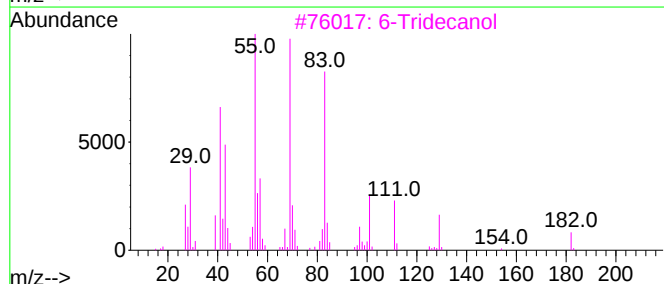
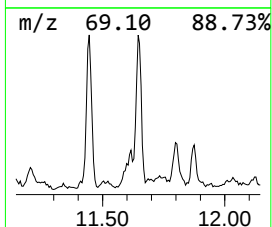
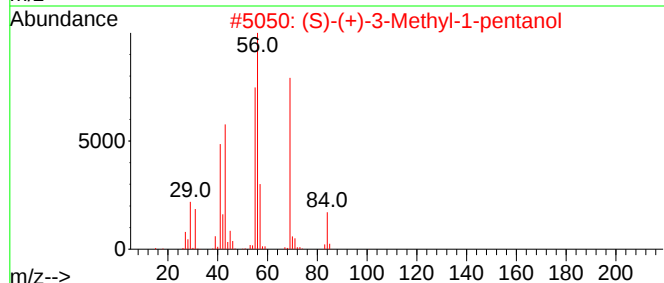
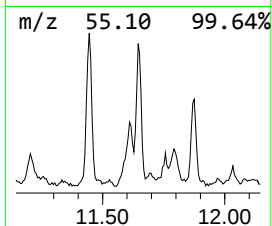
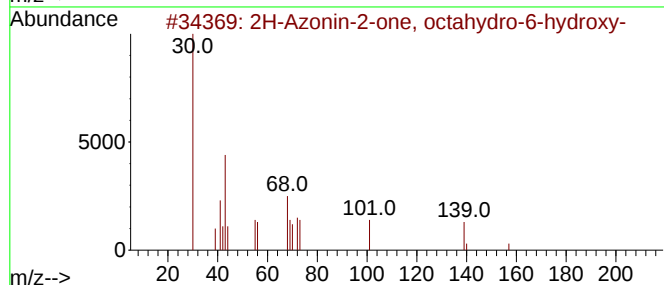
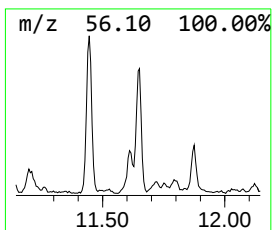
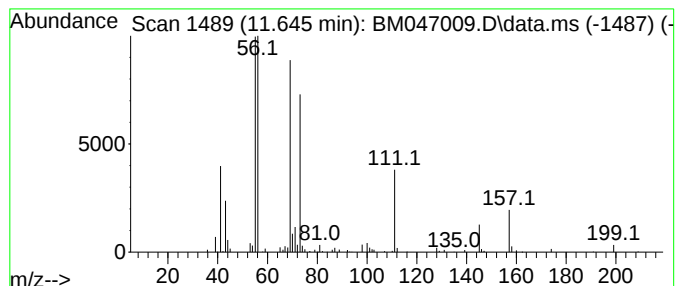
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-04 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.645	2.28 ng/ul	251227	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Azonin-2-one, octahydro-6-hyd...	157	C8H15NO2	023435-07-6	38
2			(S)-(+)-3-Methyl-1-pentanol	102	C6H14O	042072-39-9	28
3			6-Tridecanol	200	C13H28O	005770-03-6	25
4			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	14
5			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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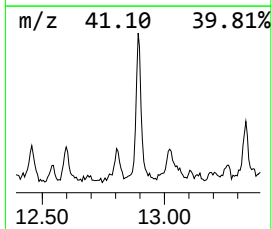
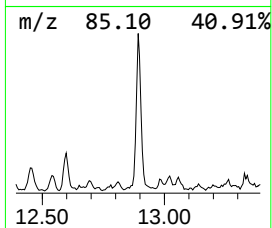
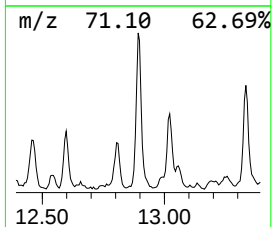
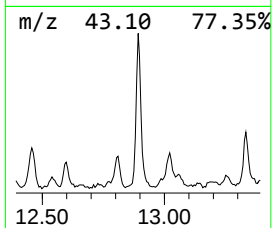
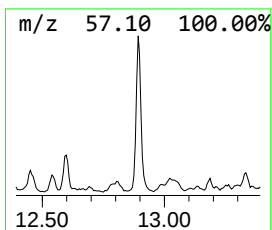
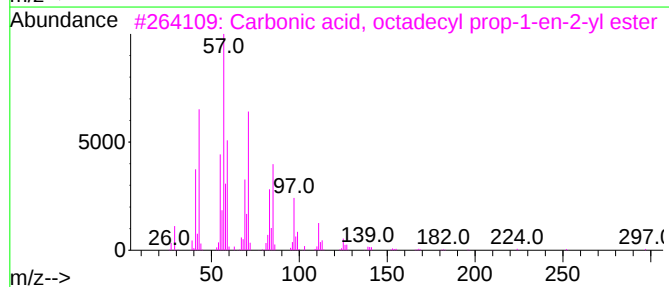
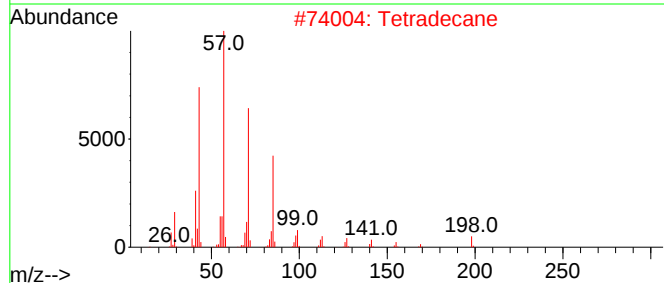
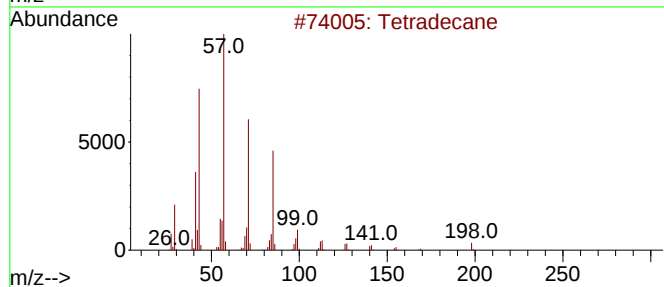
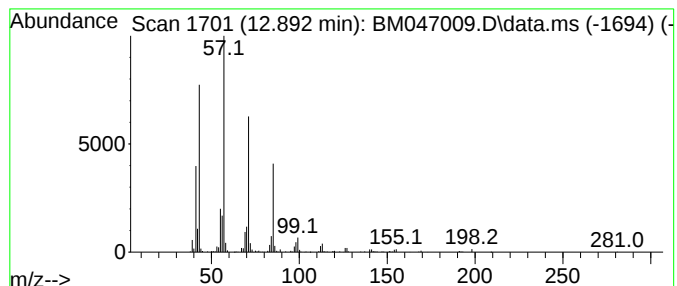
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.892	4.78 ng/ul	496862	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	96
2			Tetradecane	198	C14H30	000629-59-4	94
3			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
4			Hexadecane	226	C16H34	000544-76-3	90
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

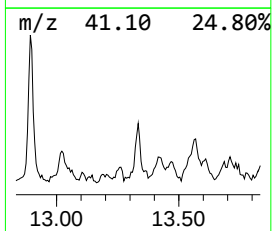
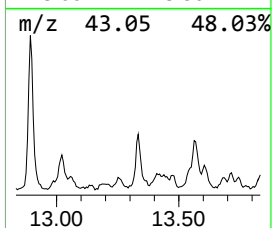
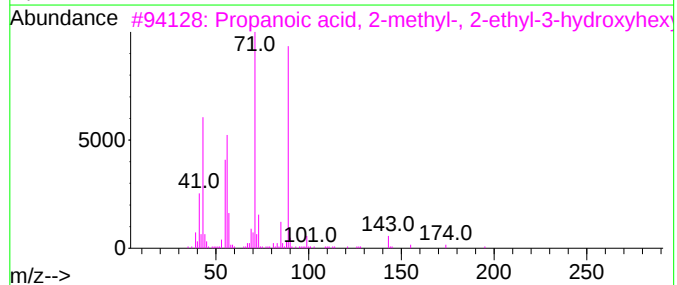
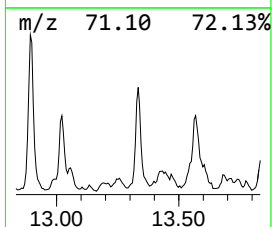
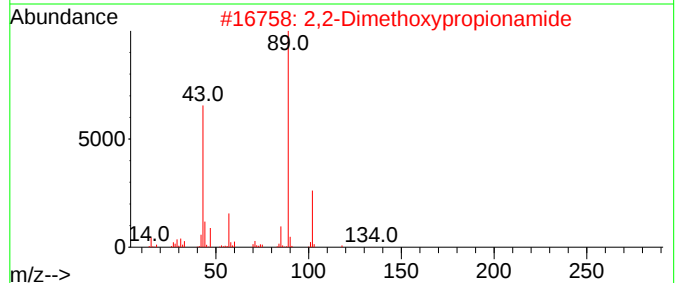
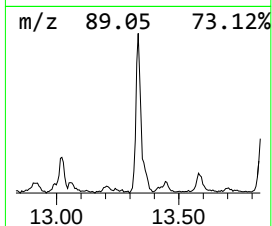
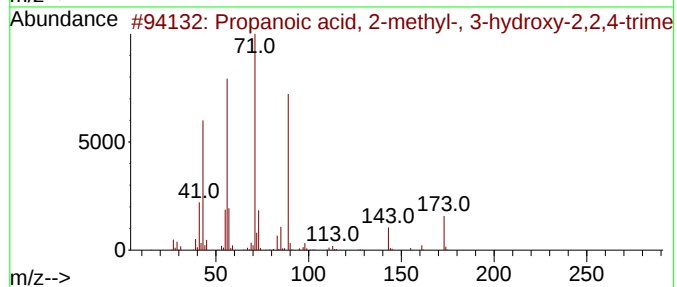
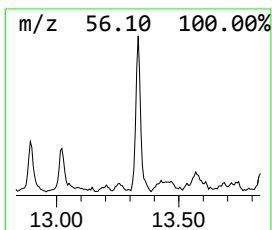
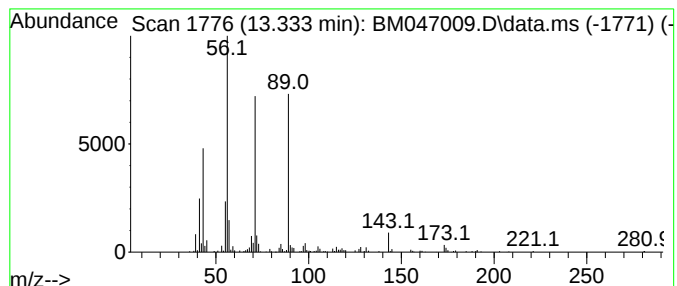
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Propanoic acid, 2-methyl-, ... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.333	2.42 ng/ul	251132	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	78
2			2,2-Dimethoxypropionamide	133	C5H11NO3	1000237-69-7	50
3			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	50
4			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	50
5			Methyl pyruvate dimethyl acetal	148	C6H12O4	010076-48-9	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId :
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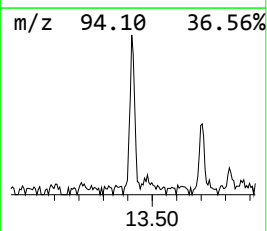
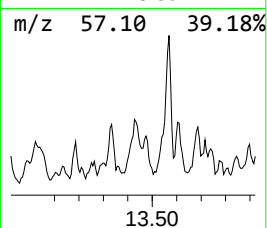
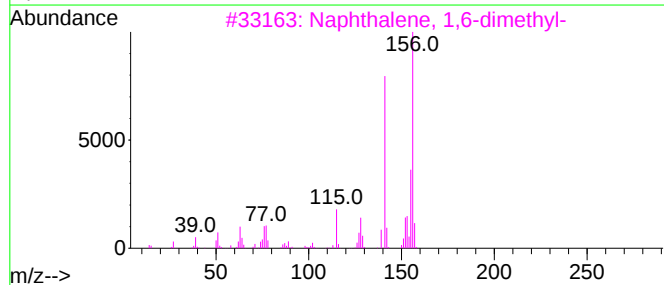
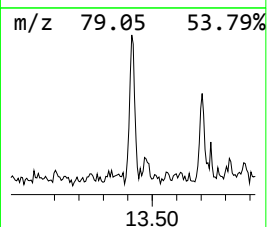
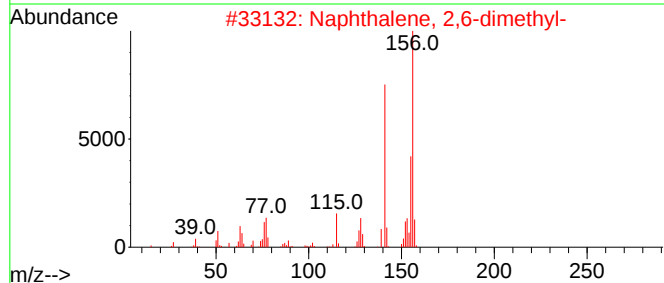
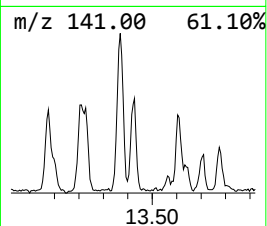
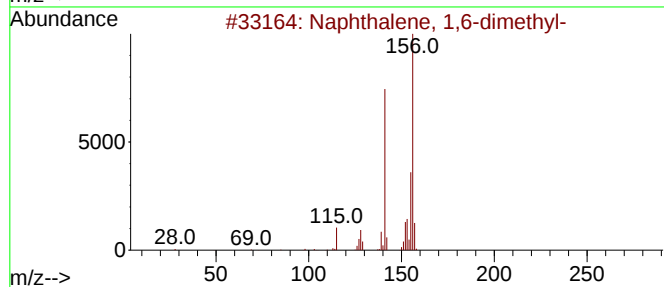
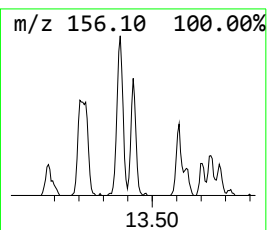
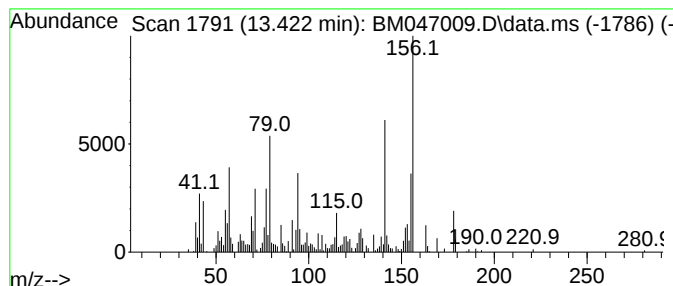
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Naphthalene, 1,6-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.422	2.09 ng/ul	217182	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

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BNA_M
ClientSampleId :
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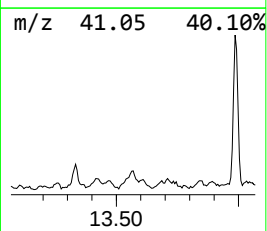
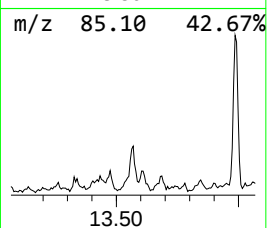
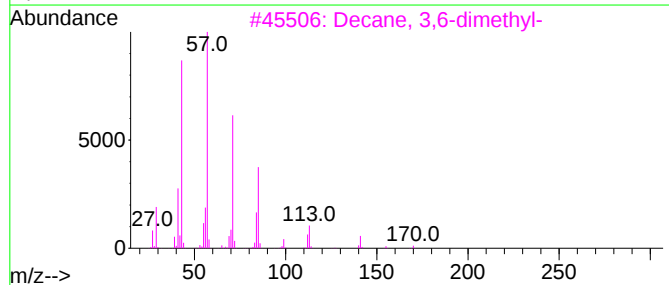
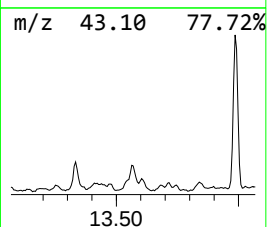
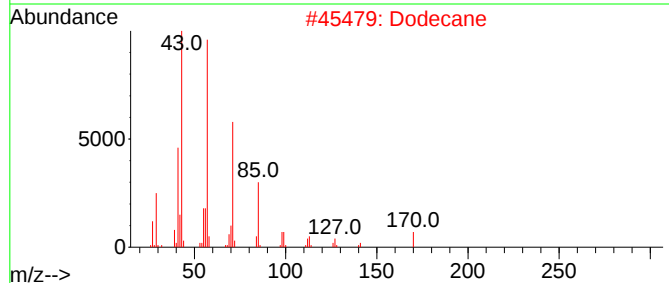
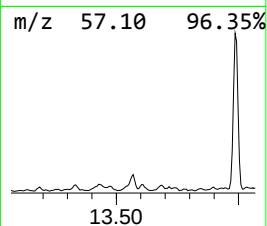
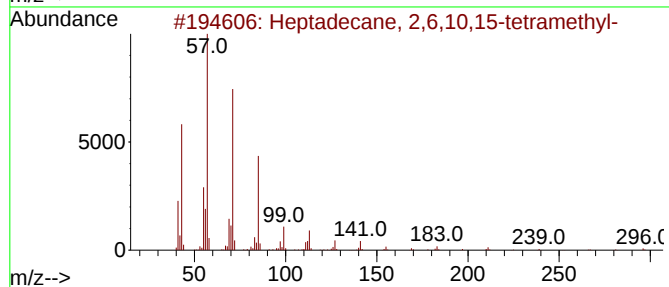
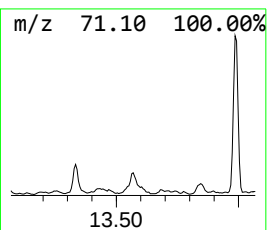
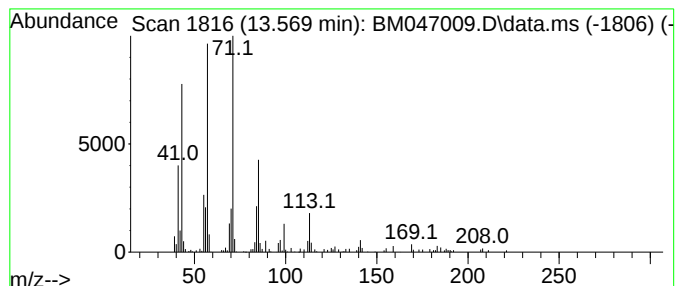
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.569	2.62 ng/ul	272554	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
2			Dodecane	170	C12H26	000112-40-3	72
3			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	52
4			Glutaric acid, butyl tetrahydrof...	272	C14H24O5	1000359-65-9	50
5			Tridecane	184	C13H28	000629-50-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

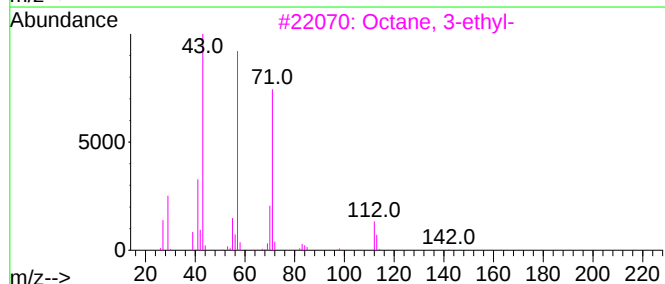
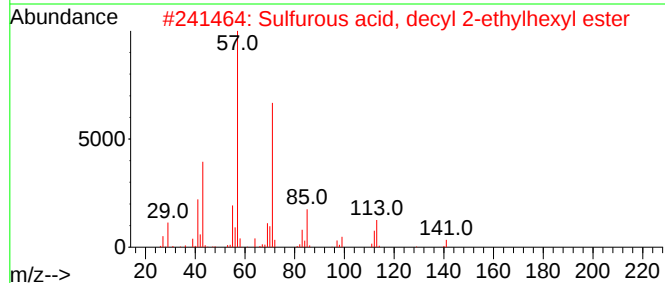
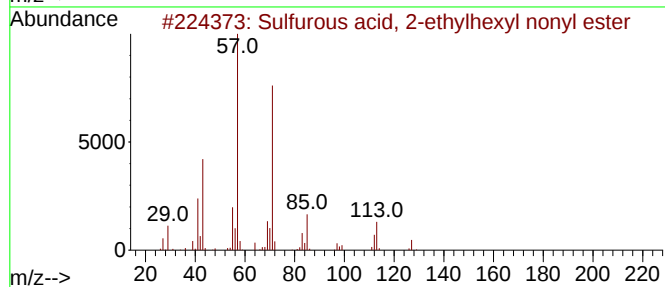
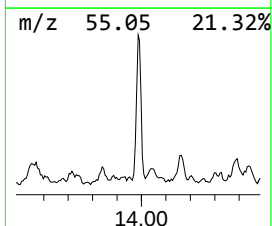
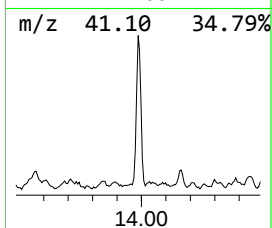
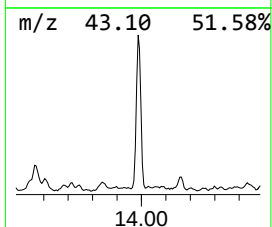
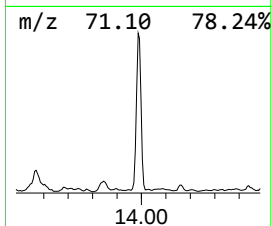
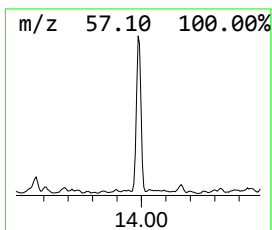
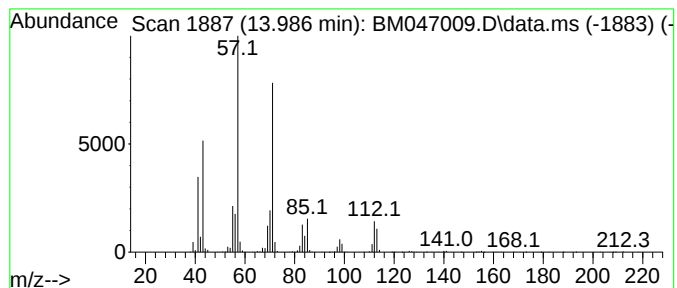
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 Sulfurous acid, 2-ethylhexy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.986	10.54 ng/ul	1095680	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72
2			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	72
3			Octane, 3-ethyl-	142	C10H22	005881-17-4	64
4			Undecane, 6-ethyl-	184	C13H28	017312-60-6	64
5			Undecane, 6-ethyl-	184	C13H28	017312-60-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

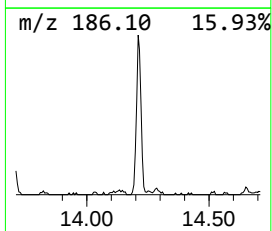
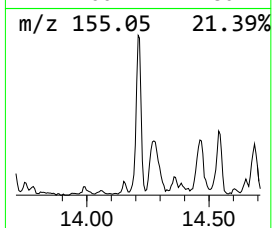
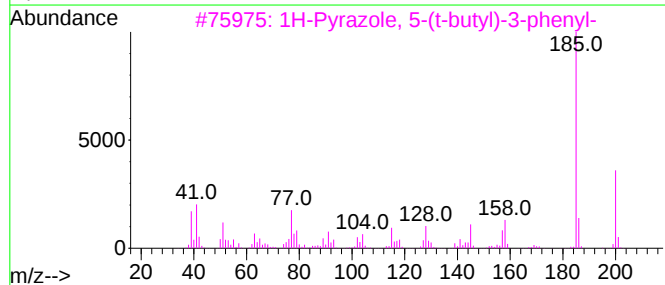
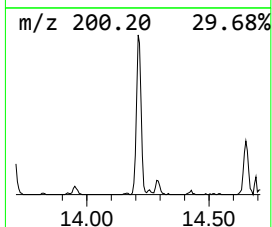
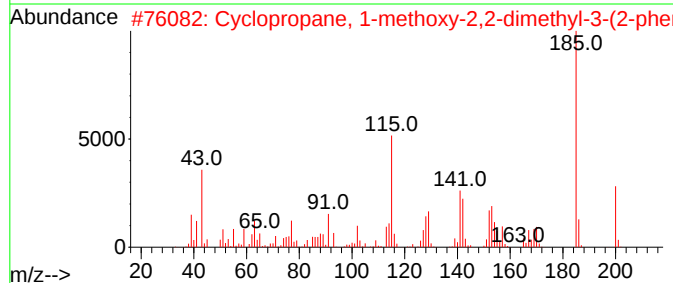
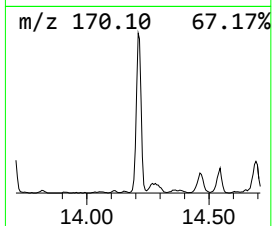
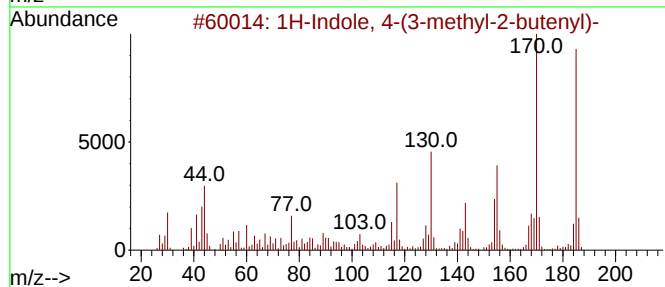
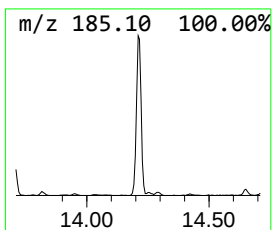
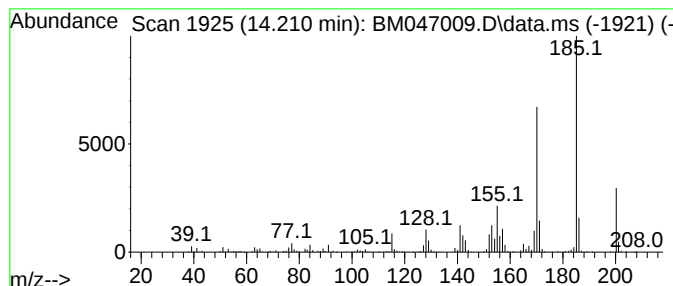
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 1H-Indole, 4-(3-methyl-2-bu... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.210	6.03 ng/ul	626386	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indole, 4-(3-methyl-2-butenyl)-	185	C13H15N	032962-25-7	74
2			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	58
3			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	55
4			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	52
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	51



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

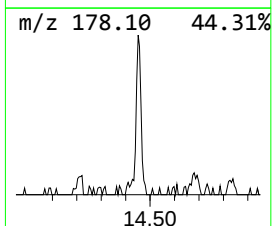
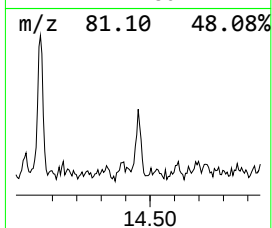
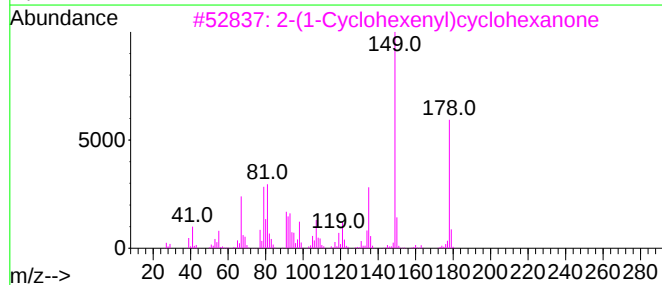
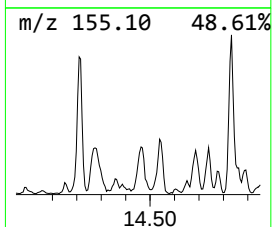
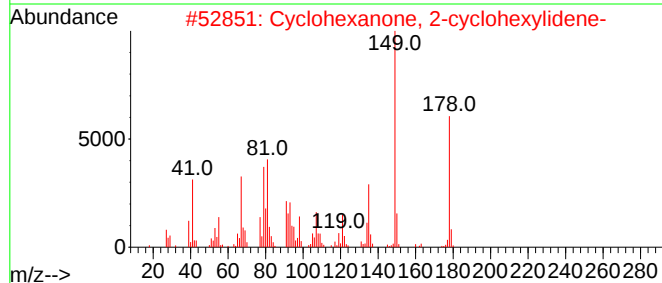
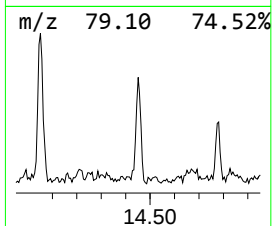
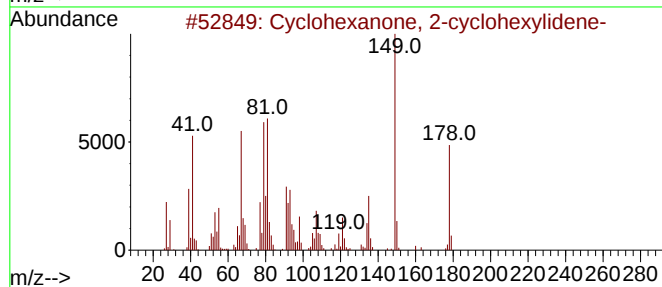
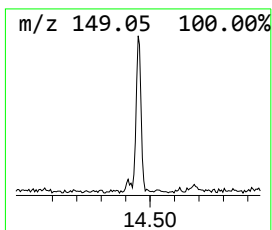
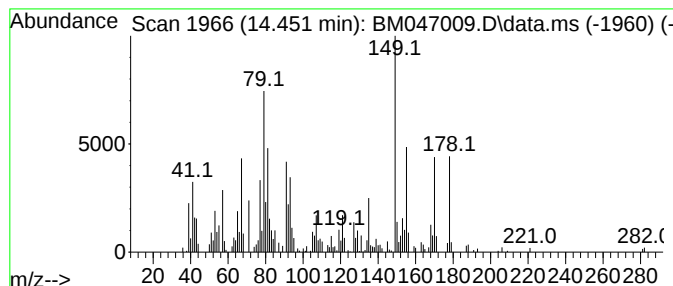
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Cyclohexanone, 2-cyclohexyl... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.451	2.38 ng/ul	247229	Acenaphthene-d10	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	83
2			Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	83
3			2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	58
4			2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	41
5			Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

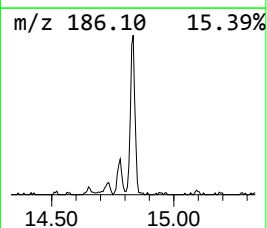
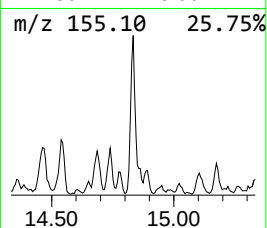
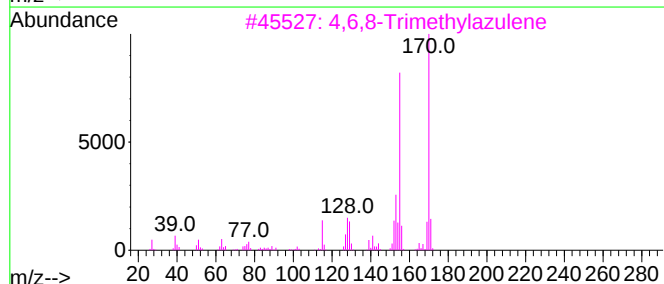
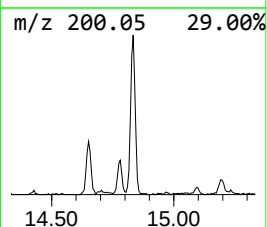
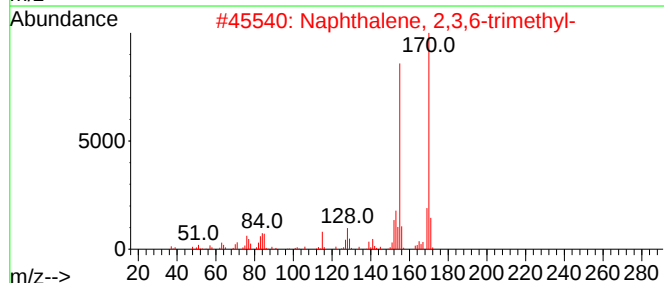
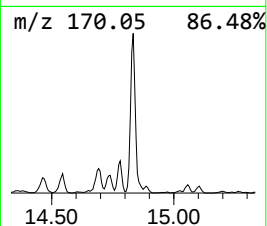
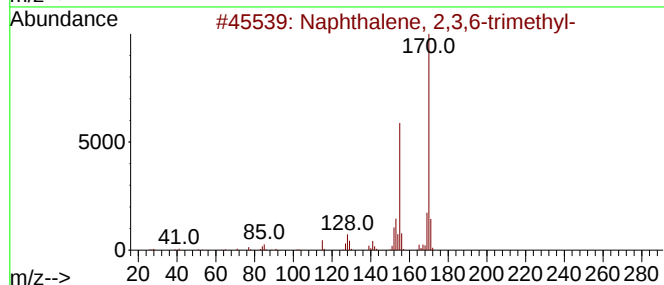
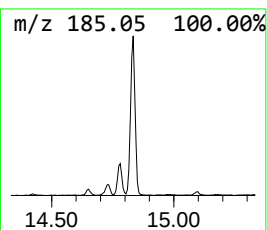
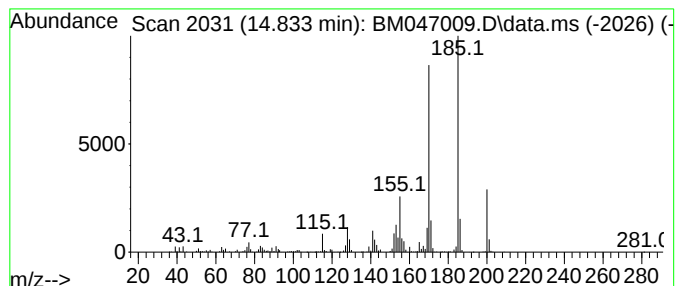
Instrument :
BNA_M
ClientSampleId :
DCVW9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 34 Naphthalene, 2,3,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.833	6.70 ng/ul	696452	Acenaphthene-d10	14.051	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	60
3	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

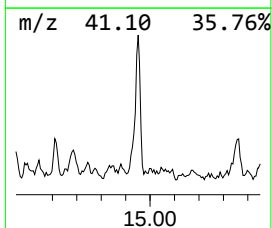
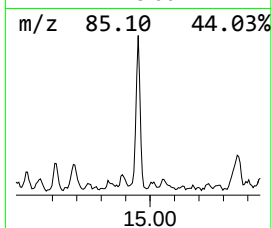
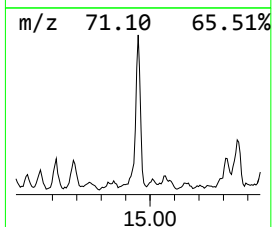
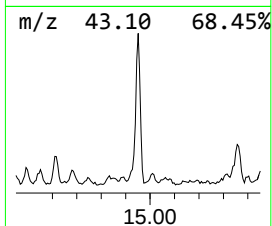
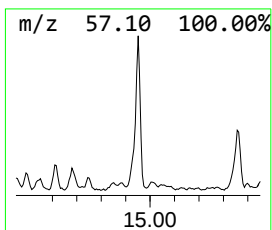
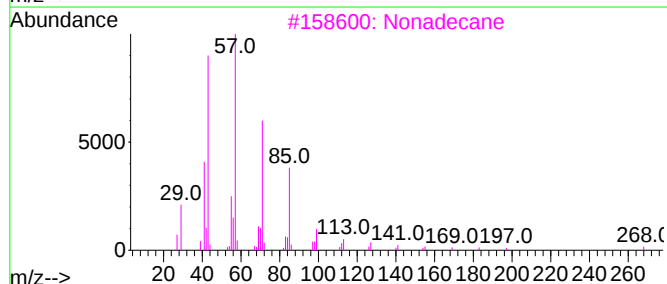
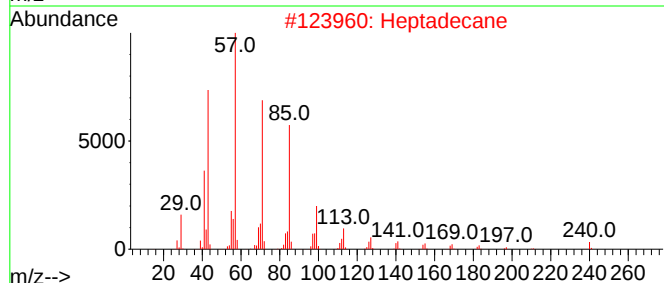
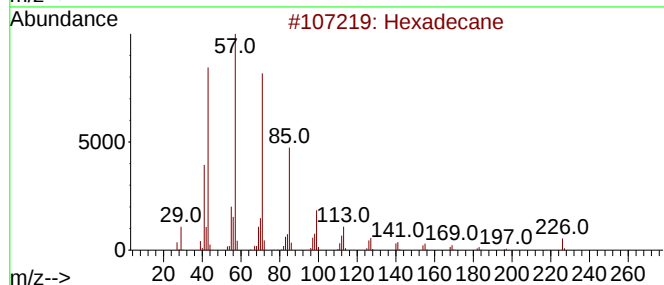
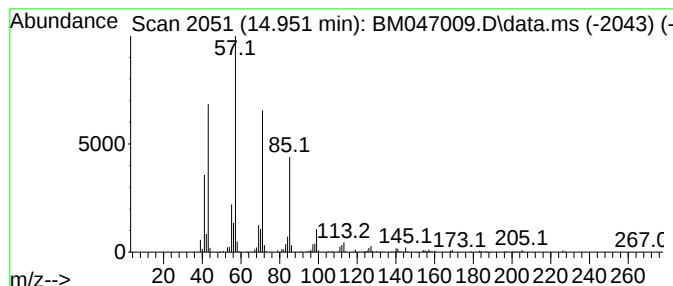
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.951	4.79 ng/ul	498290	Acenaphthene-d10	14.051

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	94
2		Heptadecane	240	C17H36	000629-78-7	90
3		Nonadecane	268	C19H40	000629-92-5	83
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	80
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

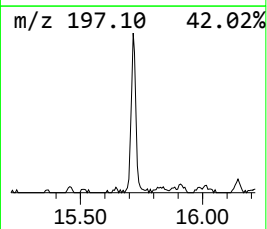
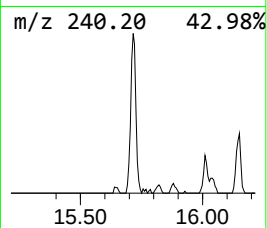
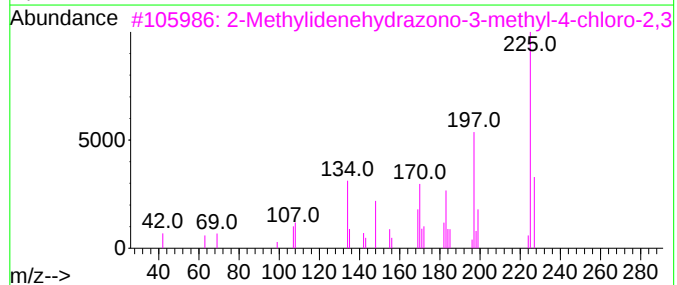
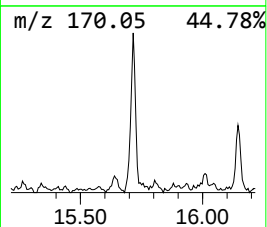
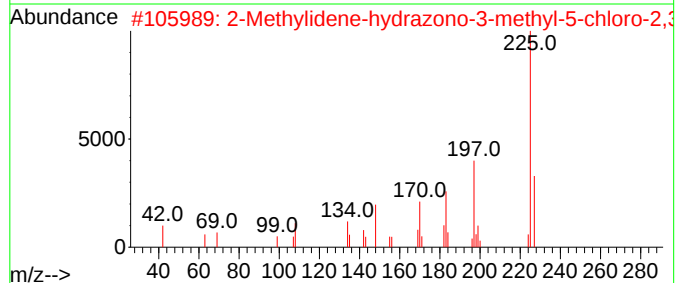
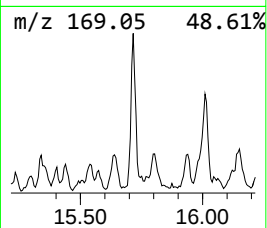
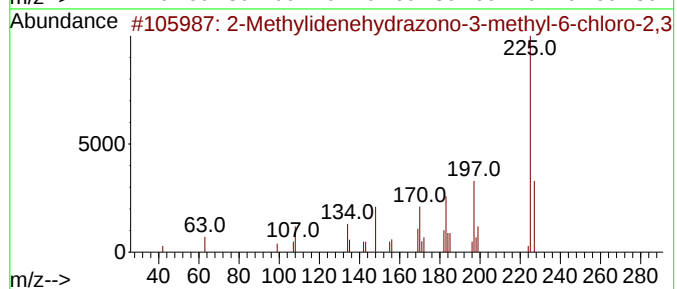
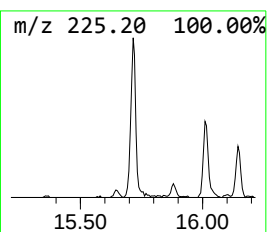
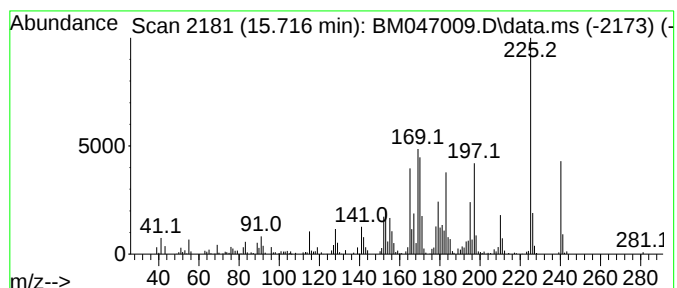
Instrument :
BNA_M
ClientSampleId :
DCVW9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 36 unknown-05 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.			
15.716	3.29 ng/ul	424355	Phenanthrene-d10	16.816			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylidenehydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-5	38		
2	2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	38		
3	2-Methylidenehydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-4	35		
4	6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	30		
5	2-Methylidenehydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-6	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

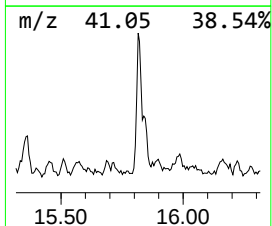
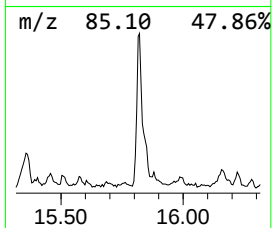
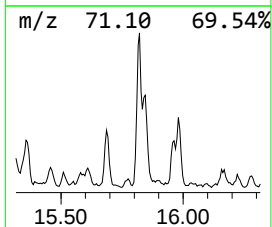
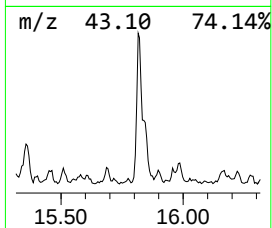
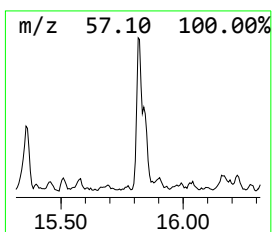
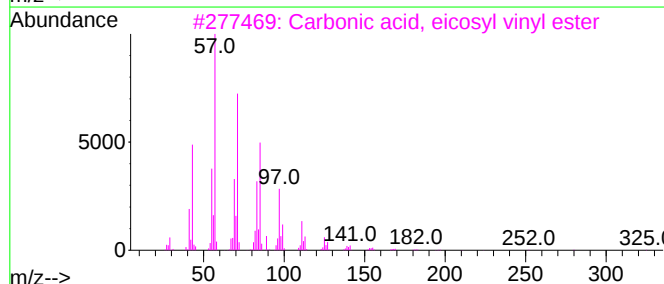
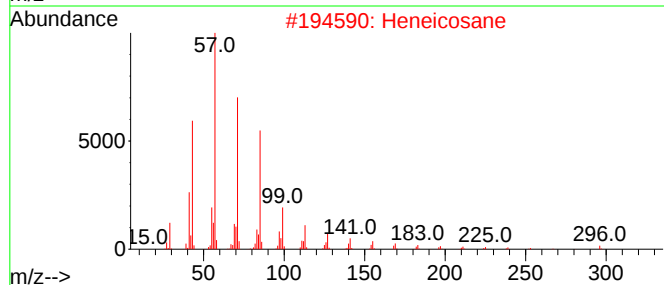
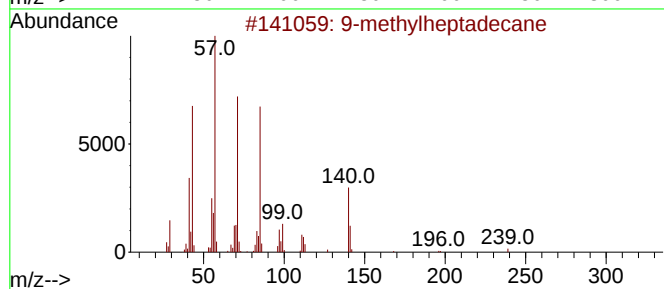
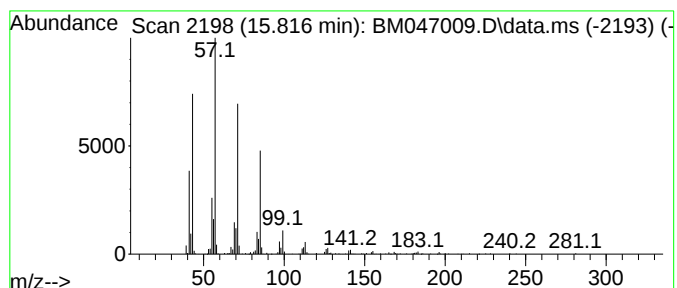
Instrument :
BNA_M
ClientSampleId :
DCVW9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.816	3.70 ng/ul	478067	Phenanthrene-d10	16.816	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-methylheptadecane	254	C18H38	026741-18-4	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86
4	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	86
5	Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

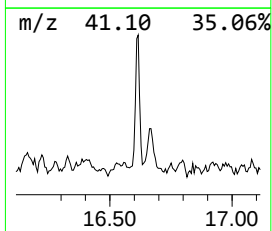
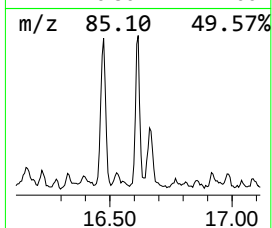
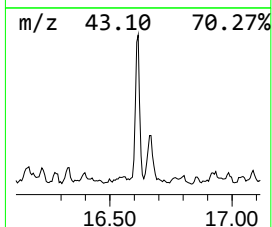
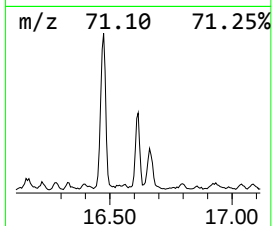
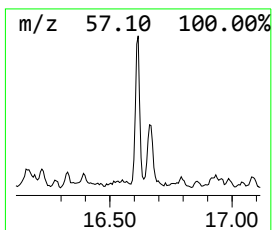
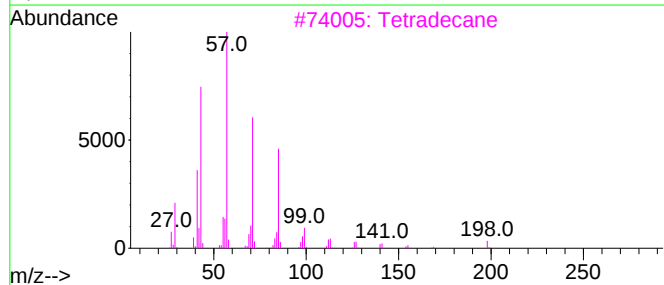
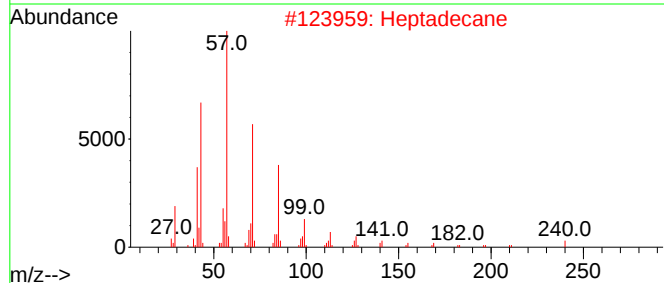
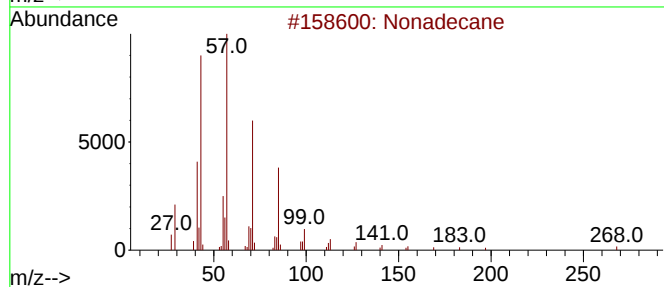
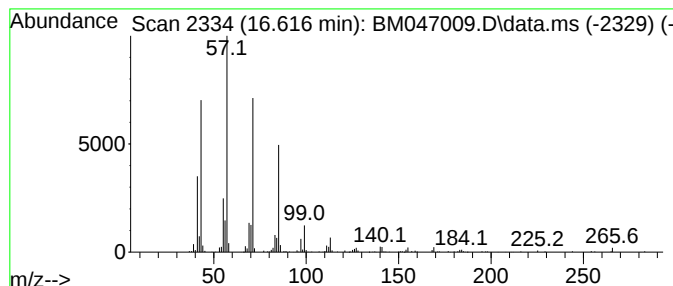
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.616	2.65 ng/ul	342671	Phenanthrene-d10	16.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Tetradecane	198	C14H30	000629-59-4	87
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	83
5		Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047009.D
Acq On : 06 Aug 2024 02:47
Operator : RC/JU
Sample : P3171-05DL 30X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCVW9DL

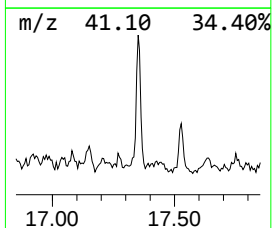
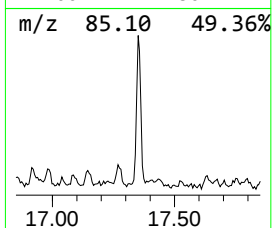
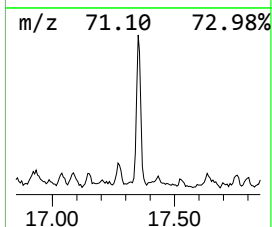
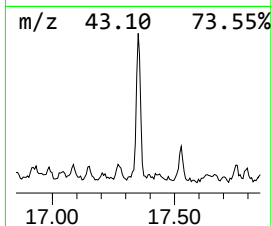
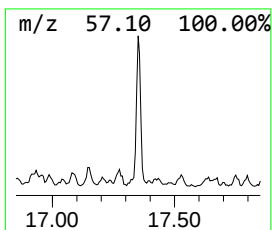
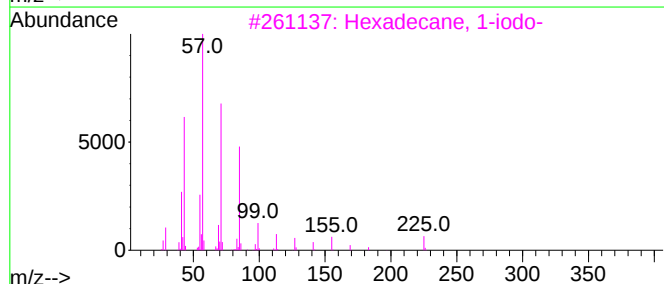
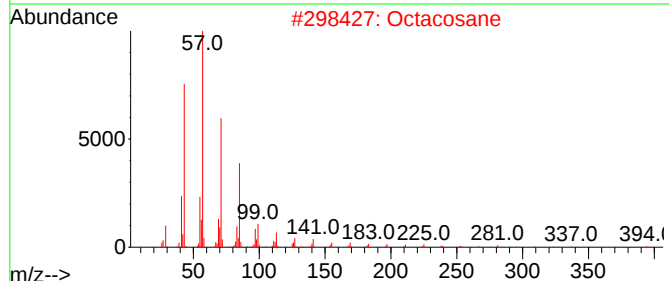
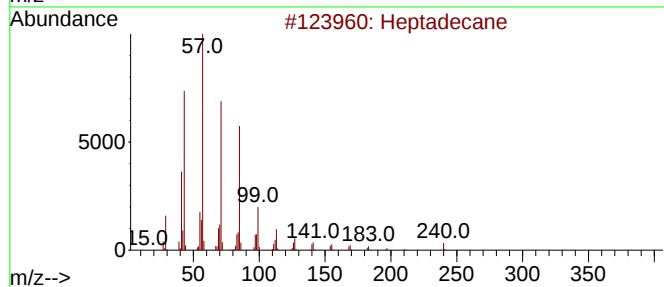
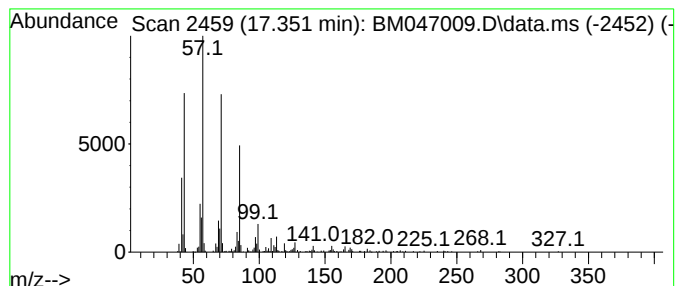
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.351	2.57 ng/ul	331661	Phenanthrene-d10	16.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	91
2		Octacosane	394	C28H58	000630-02-4	83
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
4		Heneicosane	296	C21H44	000629-94-7	80
5		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047009.D
 Acq On : 06 Aug 2024 02:47
 Operator : RC/JU
 Sample : P3171-05DL 30X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCVW9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.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...	5.463	6.3	ng/ul	377853	1	7.398	1207710	20.0
Hexylene glycol	5.775	4.4	ng/ul	264976	1	7.398	1207710	20.0
(DEL) Alkane: C...	6.063	2.2	ng/ul	133588	1	7.398	1207710	20.0
(DEL) Alkane: S...	6.416	2.5	ng/ul	147650	1	7.398	1207710	20.0
(DEL) Alkane: S...	6.469	3.1	ng/ul	187087	1	7.398	1207710	20.0
Benzene, 1-ethy...	6.504	3.1	ng/ul	184651	1	7.398	1207710	20.0
2-Heptanone, 4,...	6.846	2.4	ng/ul	142251	1	7.398	1207710	20.0
Carbonic acid, ...	7.040	20.6	ng/ul	1242900	1	7.398	1207710	20.0
1-Hexanol, 2-et...	7.475	5.8	ng/ul	353087	1	7.398	1207710	20.0
Oxalic acid, cy...	7.628	9.3	ng/ul	560356	1	7.398	1207710	20.0
Cyclohexanone, ...	7.828	2.1	ng/ul	130017	1	7.398	1207710	20.0
unknown-01	7.934	4.5	ng/ul	268914	1	7.398	1207710	20.0
(DEL) Alkane: S...	7.993	2.1	ng/ul	126638	1	7.398	1207710	20.0
Benzene, 1,4-di...	8.028	3.5	ng/ul	213831	1	7.398	1207710	20.0
(DEL) Alkane: S...	8.063	3.8	ng/ul	227787	1	7.398	1207710	20.0
Benzene, 1-meth...	8.193	2.4	ng/ul	147023	1	7.398	1207710	20.0
p-Cymene	8.393	2.1	ng/ul	125303	1	7.398	1207710	20.0
Benzene, 1-ethy...	8.487	3.9	ng/ul	234229	1	7.398	1207710	20.0
(DEL) Alkane: S...	8.622	16.6	ng/ul	1003790	1	7.398	1207710	20.0
Hexanoic acid, ...	8.716	8.3	ng/ul	503381	1	7.398	1207710	20.0
unknown-02	10.551	4.4	ng/ul	483162	2	10.151	2201460	20.0
(DEL) Alkane: S...	11.204	2.5	ng/ul	271738	2	10.151	2201460	20.0
Benzene, 4-(2-b...	11.269	4.4	ng/ul	483654	2	10.151	2201460	20.0
unknown-03	11.445	3.5	ng/ul	384599	2	10.151	2201460	20.0
(DEL) Alkane: S...	11.610	4.4	ng/ul	482791	2	10.151	2201460	20.0
unknown-04	11.645	2.3	ng/ul	251227	2	10.151	2201460	20.0
(DEL) Alkane: S...	12.892	4.8	ng/ul	496862	3	14.051	2078550	20.0
Propanoic acid,...	13.333	2.4	ng/ul	251132	3	14.051	2078550	20.0
Naphthalene, 1,...	13.422	2.1	ng/ul	217182	3	14.051	2078550	20.0
(DEL) Alkane: S...	13.569	2.6	ng/ul	272554	3	14.051	2078550	20.0
Sulfurous acid,...	13.986	10.5	ng/ul	1095680	3	14.051	2078550	20.0
1H-Indole, 4-(3...	14.210	6.0	ng/ul	626386	3	14.051	2078550	20.0
Cyclohexanone, ...	14.451	2.4	ng/ul	247229	3	14.051	2078550	20.0
Naphthalene, 2,...	14.833	6.7	ng/ul	696452	3	14.051	2078550	20.0
(DEL) Alkane: S...	14.951	4.8	ng/ul	498290	3	14.051	2078550	20.0
unknown-05	15.716	3.3	ng/ul	424355	4	16.816	2582490	20.0
(DEL) Alkane: S...	15.816	3.7	ng/ul	478067	4	16.816	2582490	20.0
(DEL) Alkane: S...	16.616	2.6	ng/ul	342671	4	16.816	2582490	20.0
(DEL) Alkane: S...	17.351	2.6	ng/ul	331661	4	16.816	2582490	20.0