

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
 Data File : BM034680.D
 Acq On : 14 Apr 2022 22:50
 Operator : CG/JU
 Sample : N2316-08
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETNQ7

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-BM040922.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034680.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.243	32	35	41	rVB	34565	44459	1.93%	0.137%
2	3.596	92	95	102	rVB	18859	31352	1.36%	0.097%
3	3.666	104	107	116	rBV	328510	501053	21.78%	1.547%
4	4.013	162	166	175	rVB3	31600	48022	2.09%	0.148%
5	4.231	200	203	210	rVB5	25263	36750	1.60%	0.113%
6	4.301	211	215	220	rVB2	12390	18571	0.81%	0.057%
7	4.672	273	278	288	rBV	80146	135973	5.91%	0.420%
8	4.843	303	307	315	rVB5	15062	26785	1.16%	0.083%
9	5.025	335	338	343	rVB3	12317	15897	0.69%	0.049%
10	5.154	355	360	366	rVB	37959	59745	2.60%	0.184%
11	5.319	383	388	400	rBV5	15890	49374	2.15%	0.152%
12	5.937	488	493	503	rVB2	33835	55341	2.41%	0.171%
13	6.125	520	525	531	rVB	19502	29977	1.30%	0.093%
14	6.537	590	595	600	rVB5	13761	24694	1.07%	0.076%
15	6.713	621	625	628	rBV3	9738	13912	0.60%	0.043%
16	6.760	630	633	637	rVB	12634	17716	0.77%	0.055%
17	6.842	642	647	650	rBV	11590	19533	0.85%	0.060%
18	7.025	672	678	698	rBV2	428022	1054803	45.85%	3.256%
19	7.190	700	706	714	rBV	362162	567124	24.65%	1.751%
20	7.389	733	740	755	rBV	456236	793295	34.48%	2.449%
21	7.866	813	821	828	rBV	452210	756087	32.87%	2.334%
22	7.937	828	833	838	rVB	38731	64681	2.81%	0.200%
23	8.160	865	871	878	rVB	86667	143479	6.24%	0.443%
24	8.331	896	900	911	rVB7	14570	30357	1.32%	0.094%
25	8.431	913	917	921	rBV5	12705	22617	0.98%	0.070%
26	8.484	921	926	934	rVB	196167	310190	13.48%	0.958%
27	8.578	935	942	950	rBV	512355	876643	38.11%	2.706%
28	9.025	1005	1018	1035	rBV	427215	884062	38.43%	2.729%
29	9.242	1050	1055	1062	rVV	35144	64827	2.82%	0.200%
30	9.472	1088	1094	1098	rBV2	12761	21336	0.93%	0.066%
31	9.566	1101	1110	1116	rBV7	16351	43959	1.91%	0.136%
32	9.678	1124	1129	1135	rBV3	31118	48930	2.13%	0.151%
33	9.754	1135	1142	1155	rVV	337198	631304	27.44%	1.949%
34	10.083	1192	1198	1205	rVB2	17420	33672	1.46%	0.104%
35	10.166	1206	1212	1218	rBV3	10530	22210	0.97%	0.069%
36	10.295	1225	1234	1246	rBV	500524	889146	38.65%	2.745%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
 Title : SVOA CALIBRATION

37	10.672	1291	1298	1314	rVB	610450	1041519	45.27%	3.215%
38	10.807	1314	1321	1331	rBV	403670	779312	33.87%	2.406%
39	10.895	1331	1336	1346	rVB4	101721	193625	8.42%	0.598%
40	11.125	1367	1375	1383	rBV6	35838	76123	3.31%	0.235%
41	11.219	1386	1391	1403	rVB3	74367	126562	5.50%	0.391%
42	11.630	1449	1461	1473	rBV	239894	530272	23.05%	1.637%
43	11.719	1473	1476	1484	rVB	52686	95499	4.15%	0.295%
44	11.819	1489	1493	1500	rVV2	31085	67824	2.95%	0.209%
45	11.883	1500	1504	1513	rVV3	24529	48560	2.11%	0.150%
46	11.983	1513	1521	1525	rVV3	6454	15181	0.66%	0.047%
47	12.036	1525	1530	1535	rVV2	24088	51305	2.23%	0.158%
48	12.101	1535	1541	1550	rVV4	109777	262411	11.41%	0.810%
49	12.183	1550	1555	1567	rVV	112708	256944	11.17%	0.793%
50	12.266	1567	1569	1574	rVV4	15158	25042	1.09%	0.077%
51	12.383	1583	1589	1598	rVB	324140	505147	21.96%	1.559%
52	12.524	1608	1613	1617	rBV6	13538	26248	1.14%	0.081%
53	12.777	1652	1656	1663	rVB9	8119	18643	0.81%	0.058%
54	12.924	1676	1681	1691	rBV4	49586	115784	5.03%	0.357%
55	13.130	1712	1716	1721	rBV8	8169	15401	0.67%	0.048%
56	13.248	1725	1736	1746	rVB6	31078	84391	3.67%	0.261%
57	13.348	1748	1753	1759	rBV	46088	67287	2.92%	0.208%
58	13.442	1765	1769	1773	rBV2	39383	57776	2.51%	0.178%
59	13.566	1783	1790	1797	rVB2	54147	93543	4.07%	0.289%
60	13.836	1828	1836	1840	rBV7	8179	12525	0.54%	0.039%
61	13.918	1844	1850	1862	rBV	907595	1333193	57.95%	4.116%
62	14.054	1868	1873	1879	rVB9	6495	15209	0.66%	0.047%
63	14.207	1888	1899	1910	rBV	1013839	1551525	67.44%	4.790%
64	14.330	1913	1920	1931	rBV4	39478	81807	3.56%	0.253%
65	14.454	1936	1941	1945	rVV3	56782	104861	4.56%	0.324%
66	14.513	1945	1951	1957	rVB	894358	1299116	56.47%	4.010%
67	14.718	1976	1986	1999	rBV	119818	368348	16.01%	1.137%
68	15.148	2053	2059	2060	rBV5	9976	13198	0.57%	0.041%
69	15.177	2060	2064	2068	rVB6	13469	20386	0.89%	0.063%
70	15.383	2093	2099	2106	rVV8	8933	23541	1.02%	0.073%
71	15.501	2112	2119	2131	rBV	1363093	1935754	84.14%	5.976%
72	15.618	2133	2139	2156	rVB	392969	603431	26.23%	1.863%
73	15.742	2156	2160	2164	rBV	13553	19576	0.85%	0.060%
74	15.865	2176	2181	2187	rVB9	6765	13501	0.59%	0.042%
75	15.930	2187	2192	2198	rVV9	7595	17386	0.76%	0.054%
76	16.177	2224	2234	2238	rBV6	16594	42898	1.86%	0.132%

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77	16.330	2255	2260	2266	rVV	81104	127483	5.54%	0.394%
78	16.395	2268	2271	2276	rVB7	9949	16269	0.71%	0.050%
79	16.483	2281	2286	2292	rBV	54726	88221	3.83%	0.272%
80	16.642	2306	2313	2318	rBV9	13430	33060	1.44%	0.102%
81	16.895	2348	2356	2359	rBV8	6804	13293	0.58%	0.041%
82	17.206	2405	2409	2411	rBV5	11034	14094	0.61%	0.044%
83	17.254	2411	2417	2427	rVV	1025866	1473677	64.06%	4.549%
84	17.354	2427	2434	2446	rVV	1333219	1948532	84.70%	6.015%
85	17.530	2459	2464	2470	rVB9	9085	18864	0.82%	0.058%
86	17.936	2528	2533	2538	rBV	25687	34568	1.50%	0.107%
87	18.153	2565	2570	2578	rBV2	160990	250202	10.88%	0.772%
88	18.948	2700	2705	2711	rVB6	11932	21309	0.93%	0.066%
89	19.083	2725	2728	2731	rBV4	9802	13197	0.57%	0.041%
90	19.165	2738	2742	2746	rBV2	54147	66162	2.88%	0.204%
91	19.212	2746	2750	2755	rVB3	19636	26411	1.15%	0.082%
92	19.289	2759	2763	2771	rBV2	17147	39767	1.73%	0.123%
93	19.430	2783	2787	2793	rBV3	32493	51195	2.23%	0.158%
94	19.642	2817	2823	2830	rBV	1638687	2300561	100.00%	7.102%
95	19.942	2872	2874	2879	rBV5	12165	18975	0.82%	0.059%
96	20.177	2909	2914	2919	rBV3	27167	46687	2.03%	0.144%
97	21.142	3074	3078	3088	rBV	466097	617127	26.83%	1.905%
98	21.430	3122	3127	3133	rBV	952935	1360586	59.14%	4.200%
99	23.647	3498	3504	3511	rBV2	926774	1920878	83.50%	5.930%
100	23.800	3524	3530	3540	rVB2	719812	1493386	64.91%	4.610%

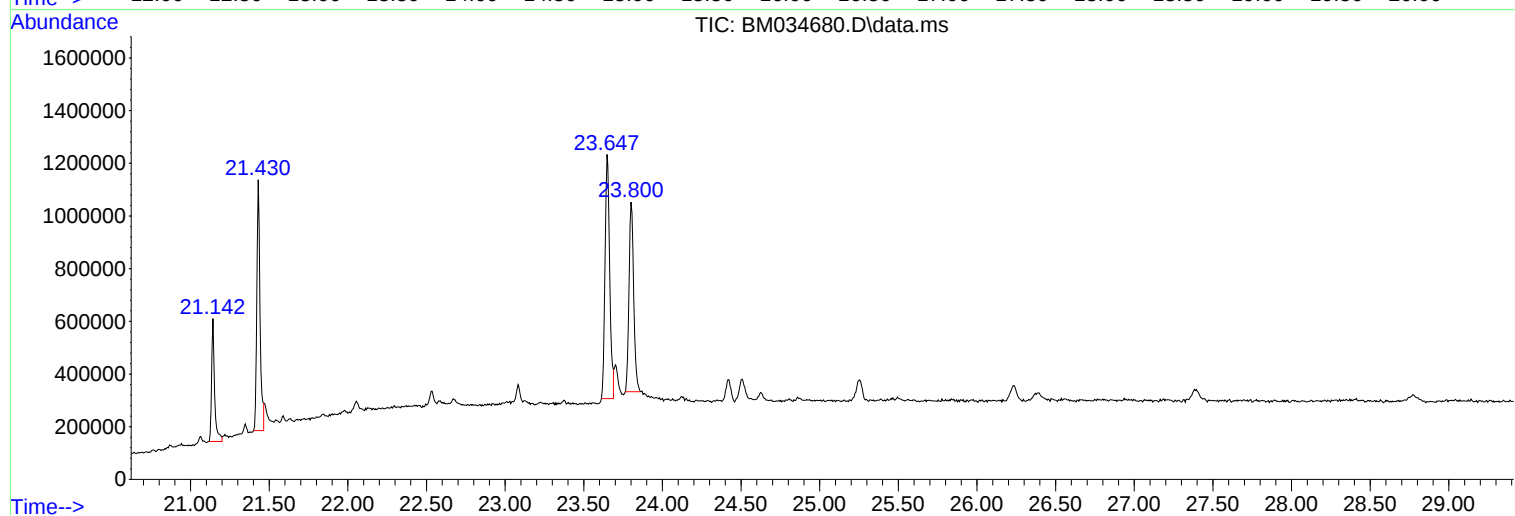
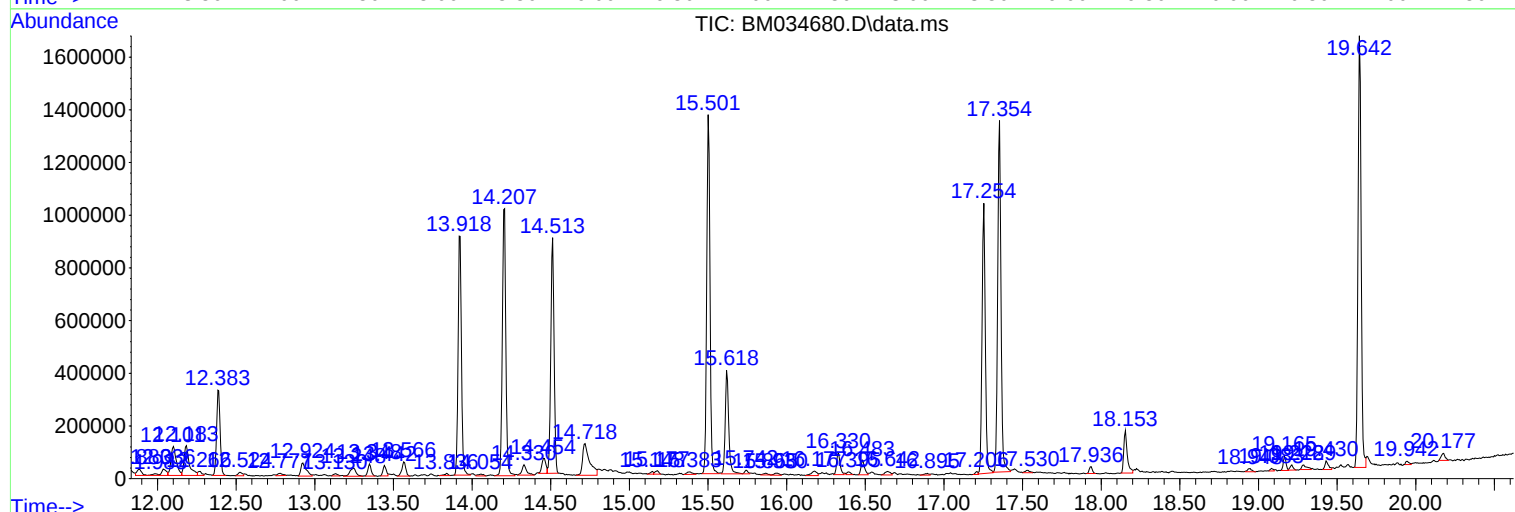
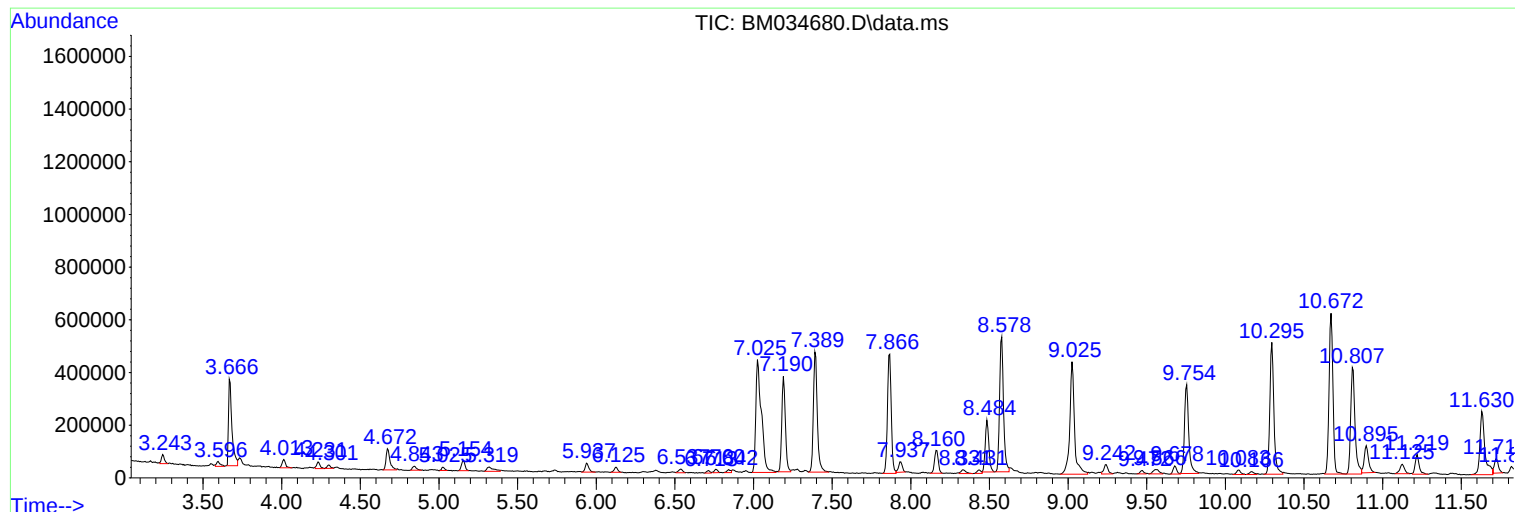
Sum of corrected areas: 32393034

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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



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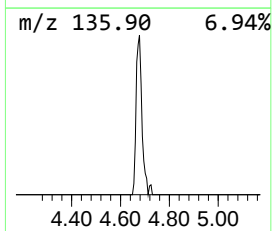
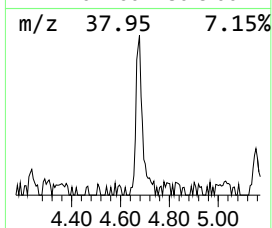
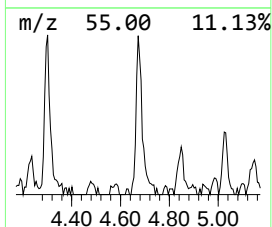
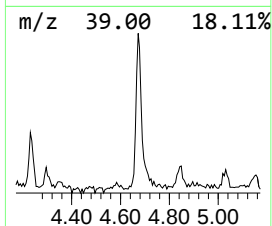
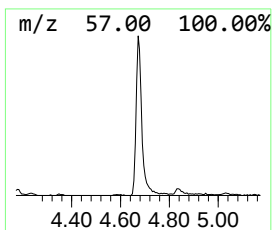
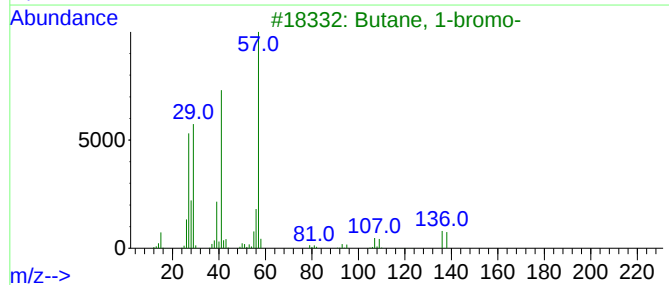
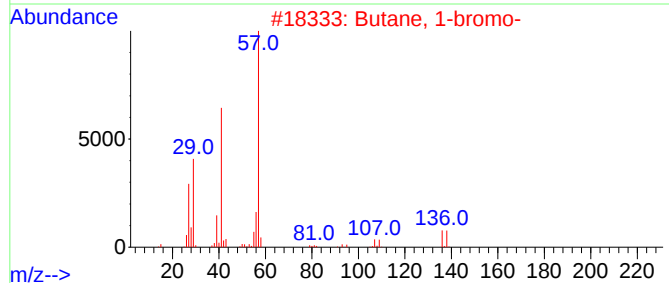
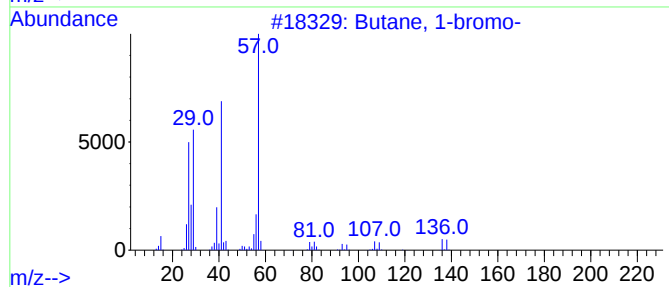
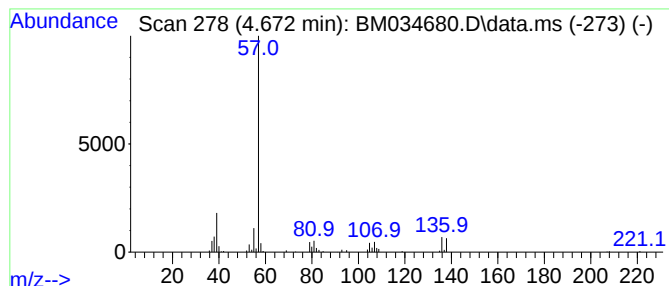
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 1-bromo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.672	3.60 ng/ul	135973	1,4-Dichlorobenzene-d4	7.866

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 1-bromo-	136	C4H9Br	000109-65-9	52
2			Butane, 1-bromo-	136	C4H9Br	000109-65-9	40
3			Butane, 1-bromo-	136	C4H9Br	000109-65-9	9
4			2,2,7,7-Tetramethyl-4,5-dimethyl...	194	C14H26	090822-63-2	9
5			2-Propyn-1-ol, propionate	112	C6H8O2	001932-92-9	9



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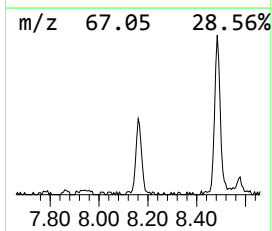
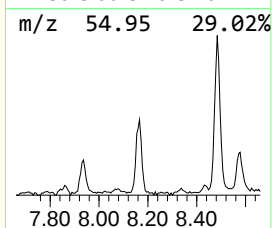
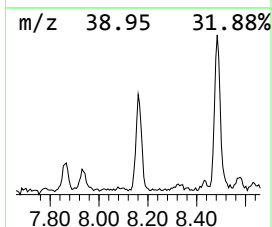
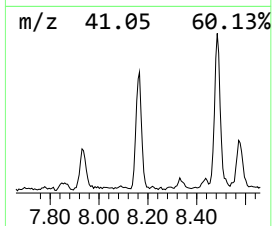
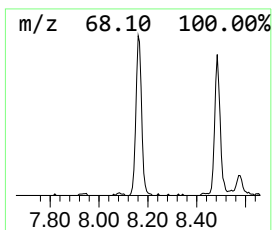
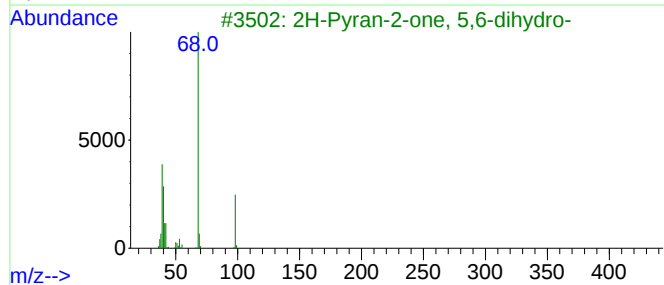
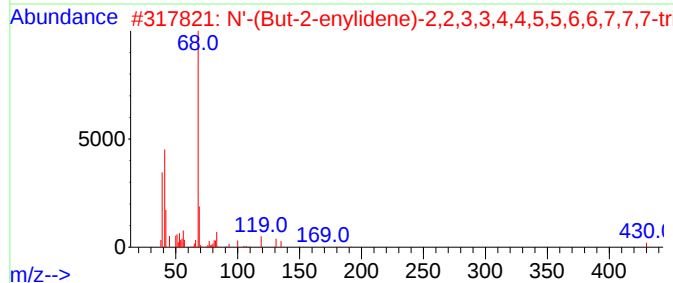
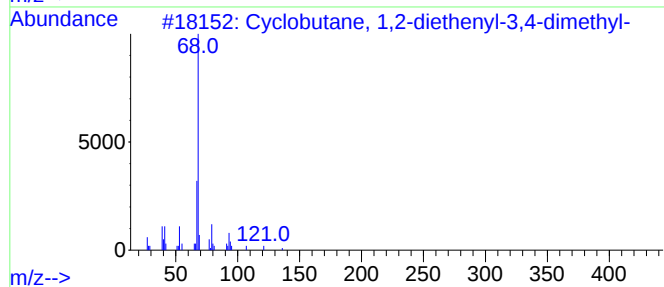
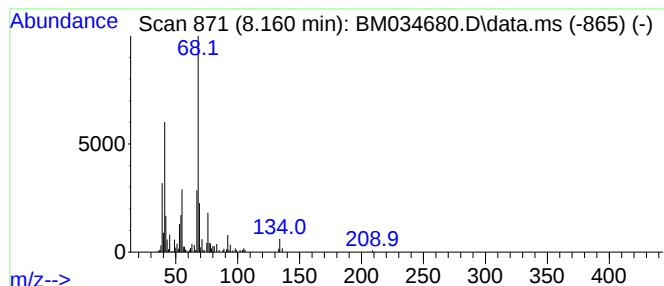
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclobutane, 1,2-diethenyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.160	3.80 ng/ul	143479	1,4-Dichlorobenzene-d4	7.866

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutane, 1,2-diethenyl-3,4-d...	136	C10H16	1000034-00-1	53
2			N'-(But-2-enylidene)-2,2,3,3,4,4...	430	C11H7F13N2O	330591-15-6	38
3			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	38
4			Cyclobutane, 1,2-dipropenyl-	136	C10H16	022769-00-2	36
5			2-Pentyne	68	C5H8	000627-21-4	36



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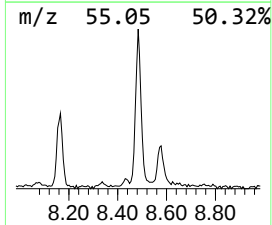
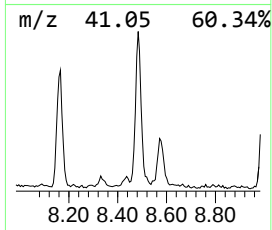
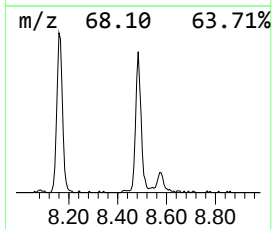
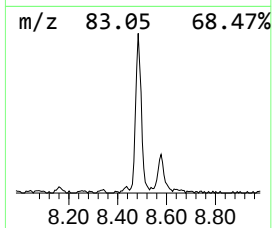
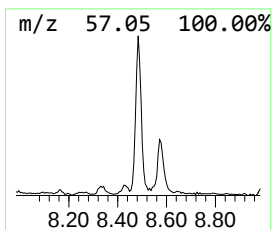
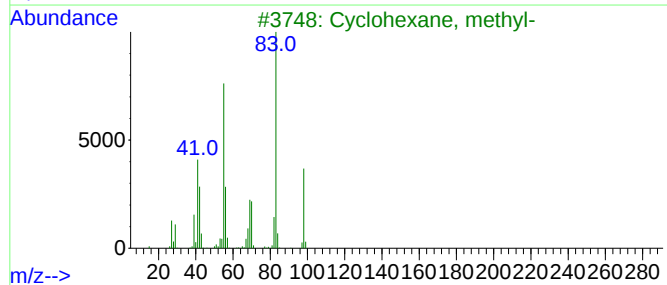
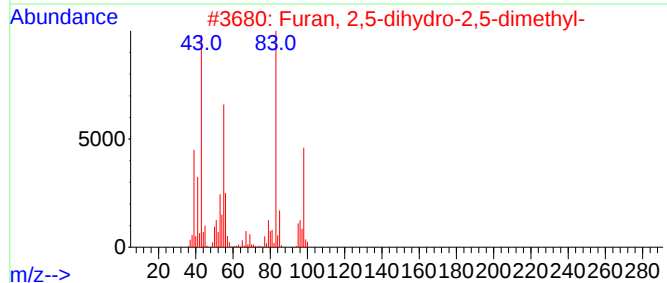
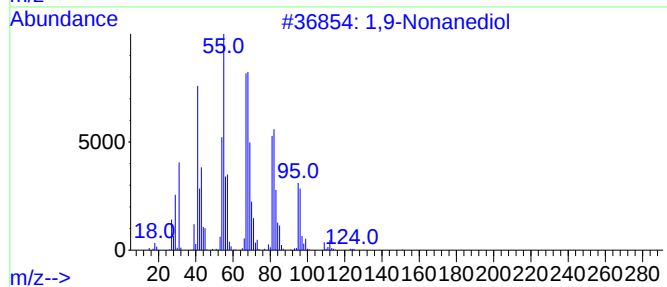
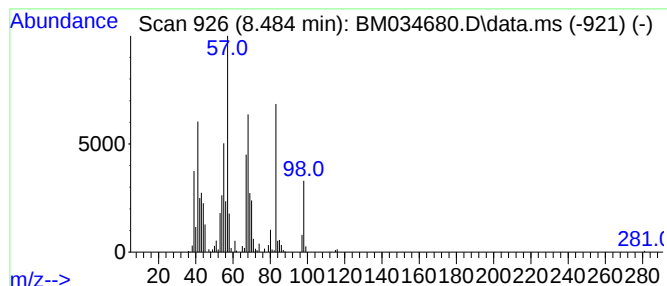
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.484	8.21 ng/ul	310190	1,4-Dichlorobenzene-d4	7.866

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,9-Nonanediol	160	C9H20O2	003937-56-2	43
2			Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	42
3			Cyclohexane, methyl-	98	C7H14	000108-87-2	30
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	30
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
Operator : CG/JU
Sample : N2316-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

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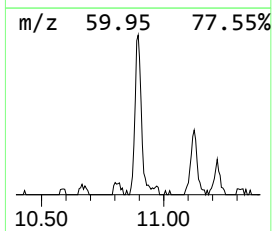
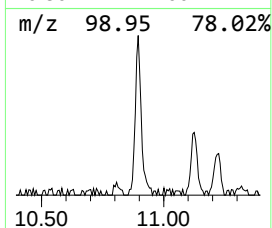
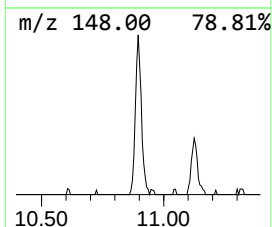
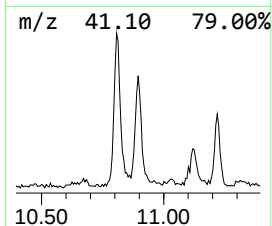
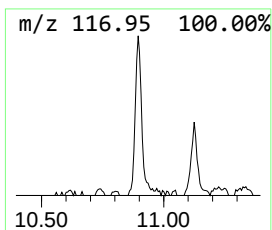
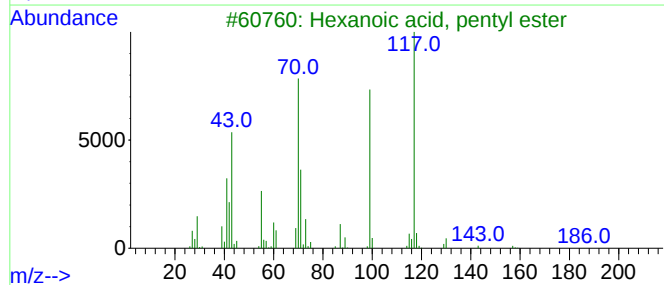
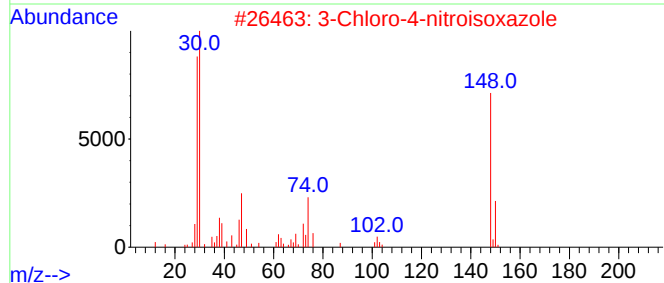
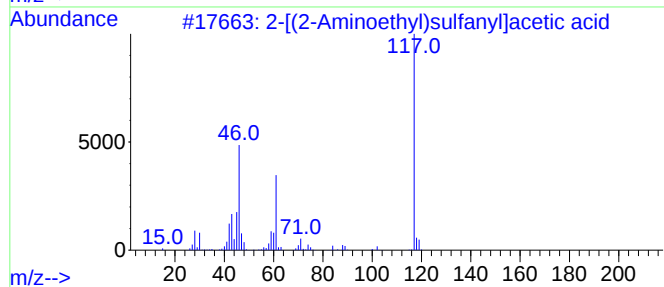
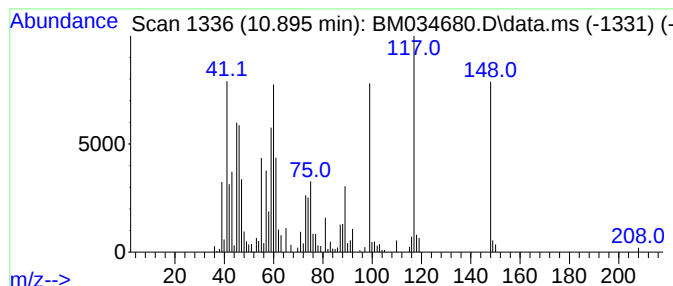
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.895	3.72 ng/ul	193625	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[(2-Aminoethyl)sulfanyl]acetic...	135	C4H9N02S	003335-52-2	22		
2	3-Chloro-4-nitroisoxazole	148	C3HC1N2O3	1000432-93-0	14		
3	Hexanoic acid, pentyl ester	186	C11H22O2	000540-07-8	12		
4	Thiirane	60	C2H4S	000420-12-2	11		
5	2,4-Thiazolidinedione	117	C3H3N02S	002295-31-0	11		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
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ClientSampleId :
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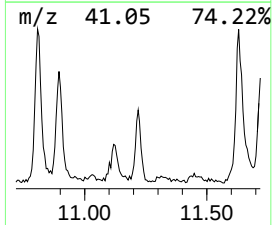
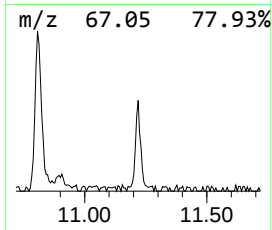
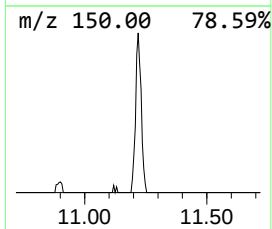
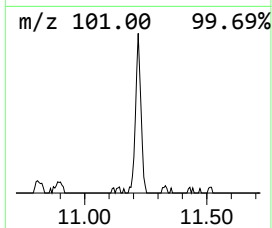
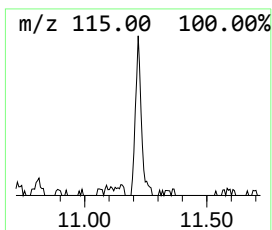
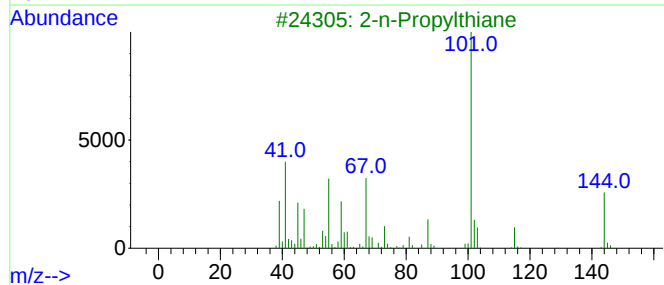
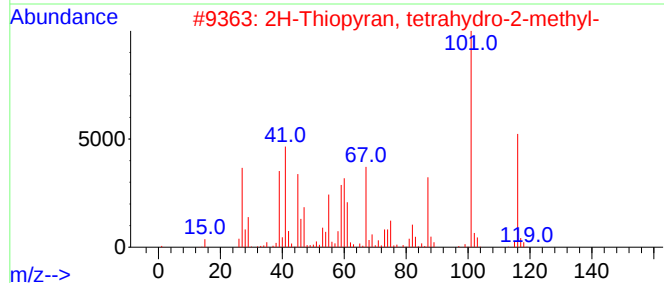
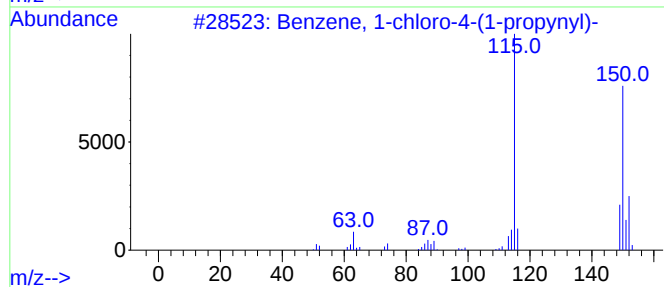
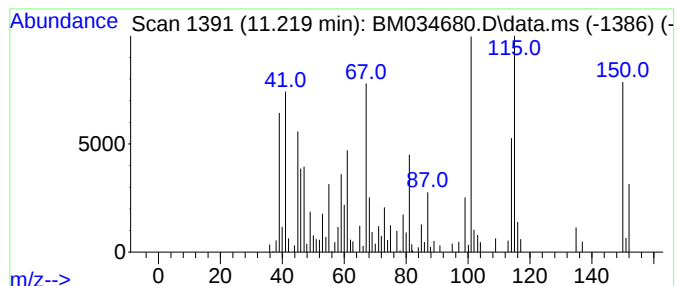
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.219	2.43 ng/ul	126562	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-(1-propynyl)-	150	C9H7Cl	002809-65-6	27
2			2H-Thiopyran, tetrahydro-2-methyl-	116	C6H12S	005161-16-0	25
3			2-n-Propylthiane	144	C8H16S	017912-23-1	22
4			2-Chloro-4,6-difluoro-pyrimidine	150	C4HClF2N2	038953-30-9	22
5			9-Chloro-bicyclo[6.1.0]nona-1(8)...	150	C9H7Cl	1010295-96-4	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
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ClientSampleId :
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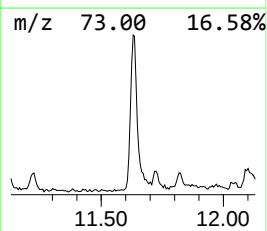
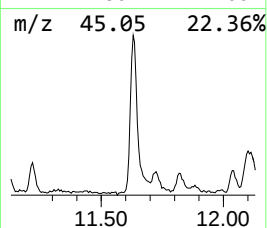
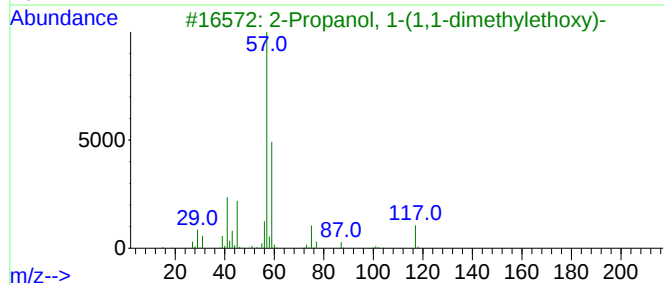
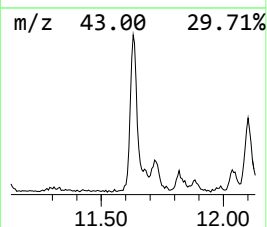
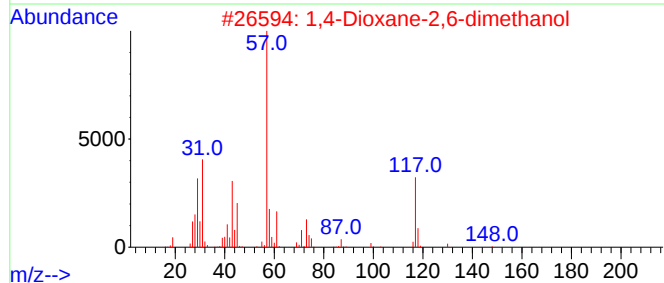
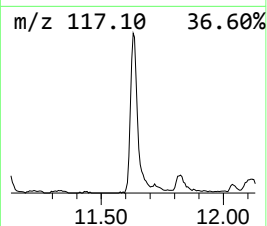
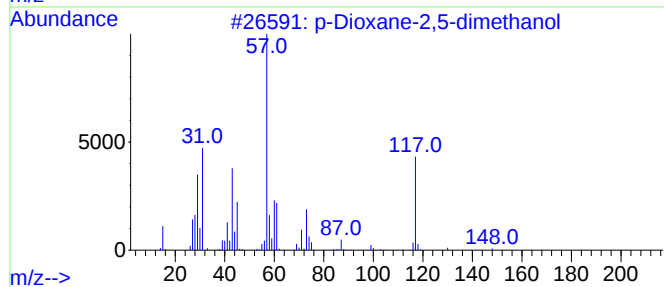
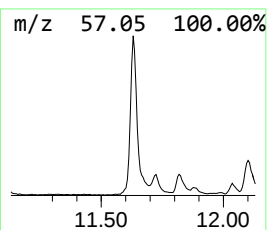
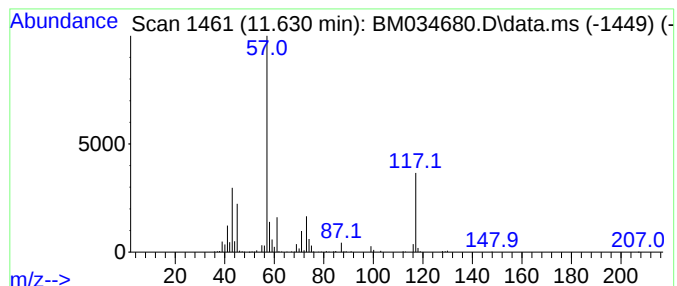
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 p-Dioxane-2,5-dimethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.630	10.18 ng/ul	530272	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Dioxane-2,5-dimethanol	148	C6H12O4	014236-12-5	83
2			1,4-Dioxane-2,6-dimethanol	148	C6H12O4	054120-69-3	45
3			2-Propanol, 1-(1,1-dimethylethoxy)-	132	C7H16O2	057018-52-7	38
4			Propanoic acid, 2,2-dimethyl-, m...	116	C6H12O2	000598-98-1	30
5			Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
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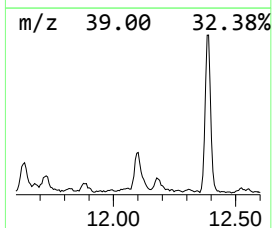
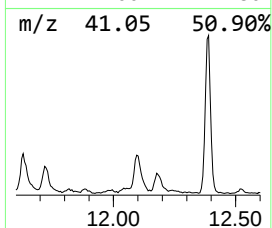
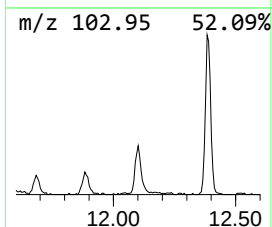
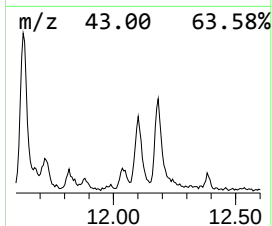
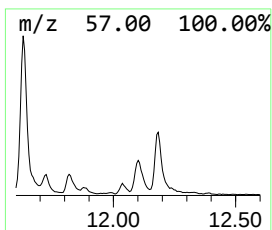
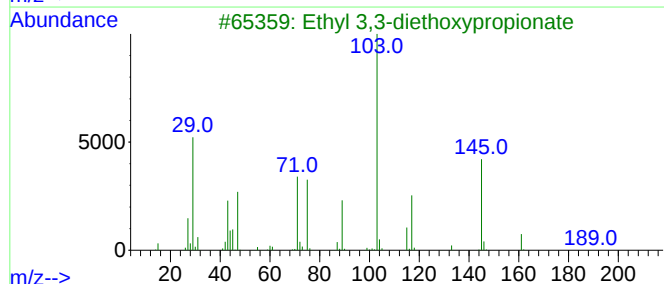
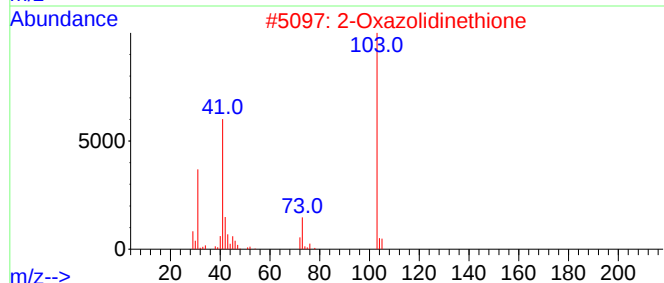
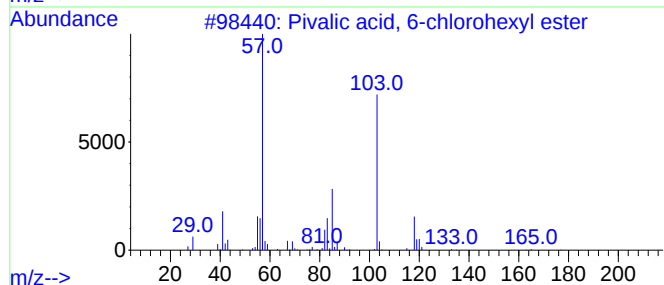
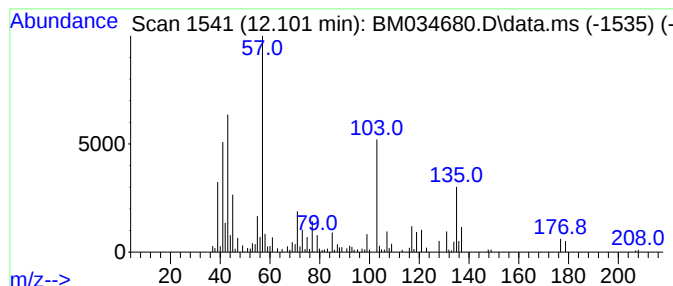
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.101	5.04 ng/ul	262411	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pivalic acid, 6-chlorohexyl ester	220	C11H21ClO2	1000330-91-8	22
2			2-Oxazolidinethione	103	C3H5NOS	005840-81-3	22
3			Ethyl 3,3-diethoxypropionate	190	C9H18O4	010601-80-6	22
4			2-t-Butyl-6-chloromethyl-[1,3]di...	206	C9H15ClO3	139883-58-2	12
5			Heptane, 1,1-diethoxy-	188	C11H24O2	000688-82-4	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
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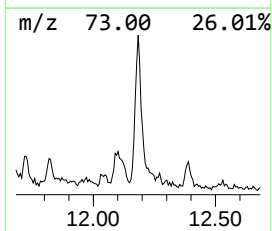
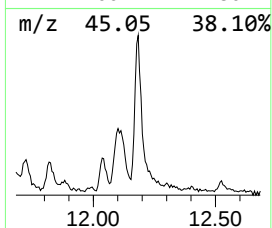
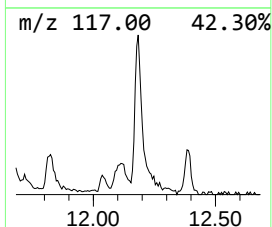
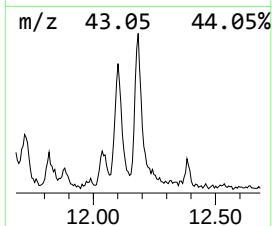
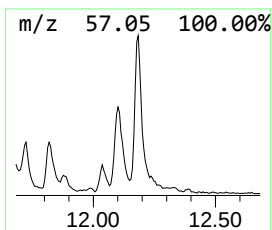
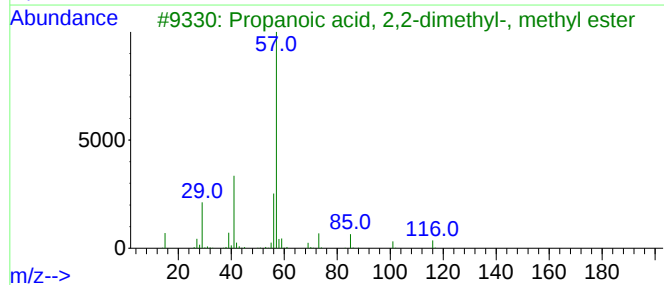
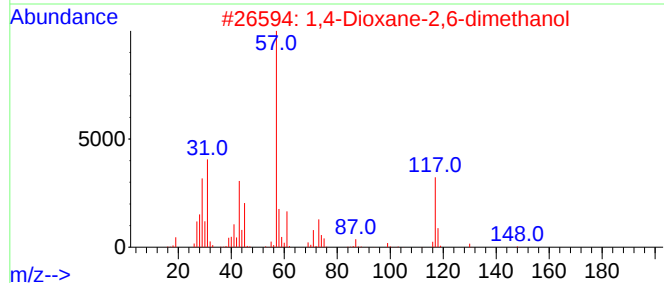
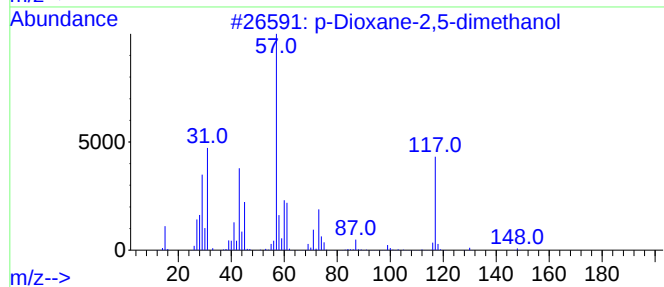
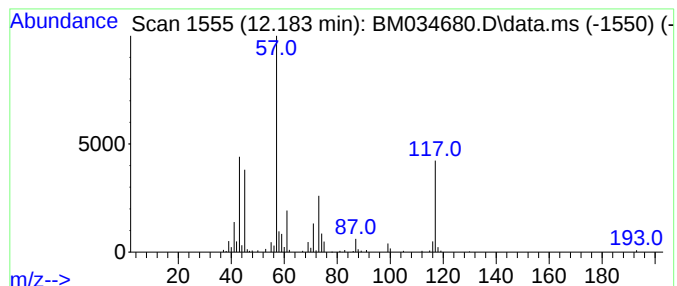
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.183	4.93 ng/ul	256944	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Dioxane-2,5-dimethanol	148	C6H12O4	014236-12-5	33
2			1,4-Dioxane-2,6-dimethanol	148	C6H12O4	054120-69-3	33
3			Propanoic acid, 2,2-dimethyl-, m...	116	C6H12O2	000598-98-1	11
4			2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	10
5			Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
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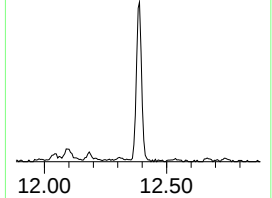
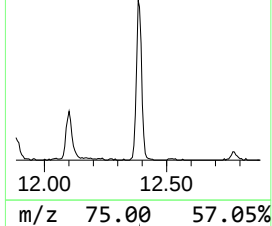
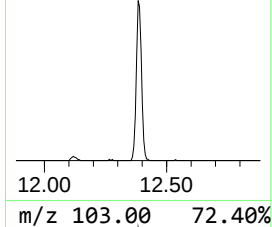
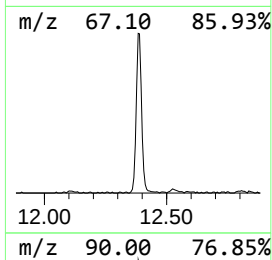
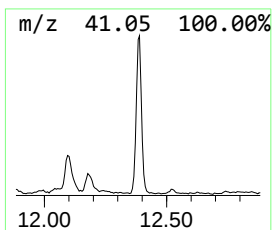
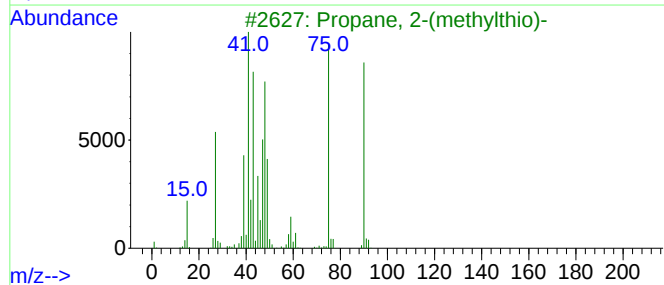
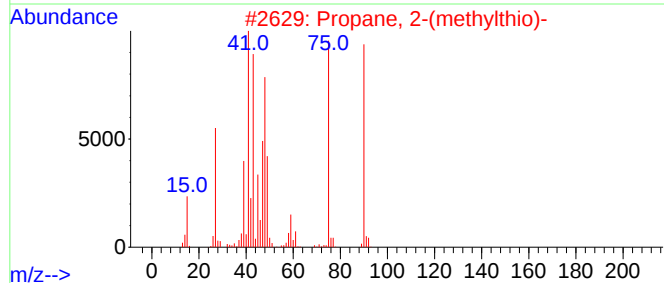
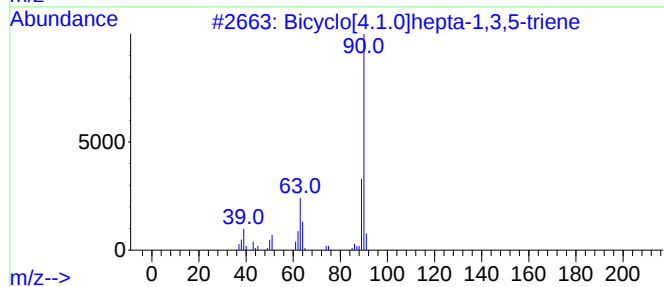
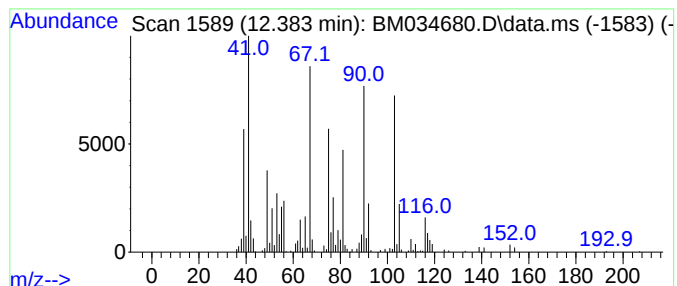
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.383	9.70 ng/ul	505147	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	43
2			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	40
3			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	40
4			1-Hexene, 3-chloro-	118	C6H11Cl	053101-38-5	37
5			Thiourea, methyl-	90	C2H6N2S	000598-52-7	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
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Sample : N2316-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

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ClientSampleId :
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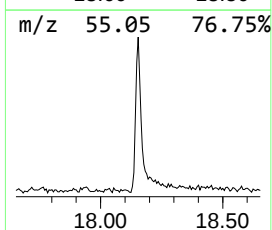
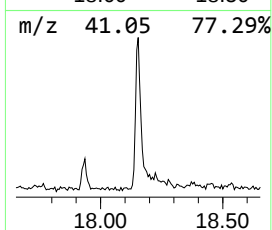
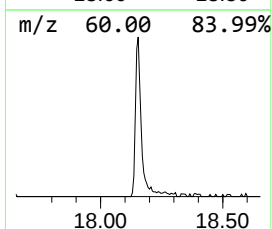
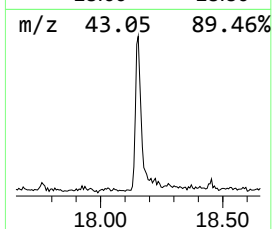
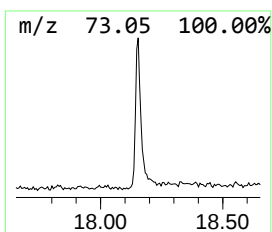
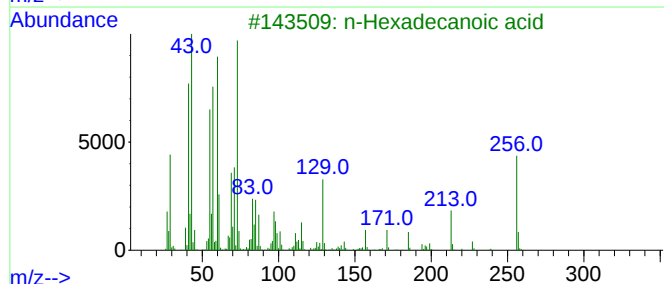
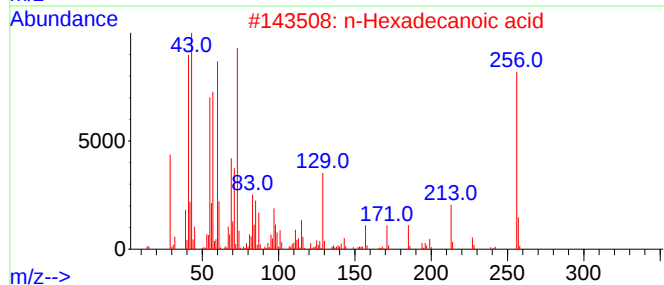
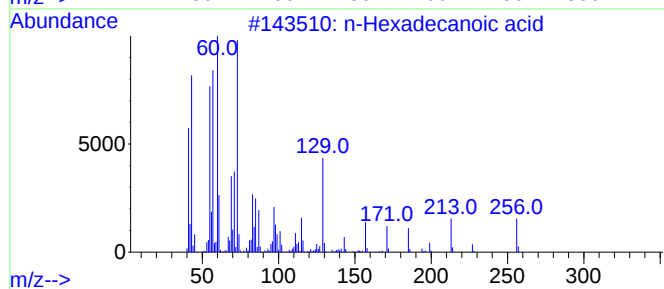
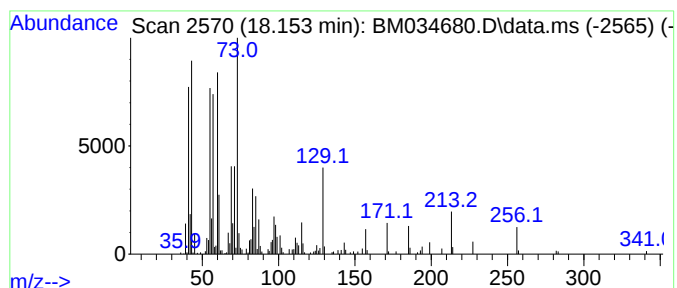
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.154	3.40 ng/ul	250202	Phenanthrene-d10	17.254

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5		Octadecanoic acid	284	C18H36O2	000057-11-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
Operator : CG/JU
Sample : N2316-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ7

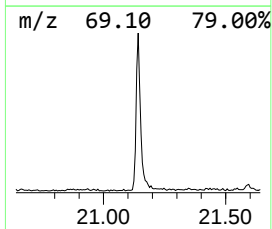
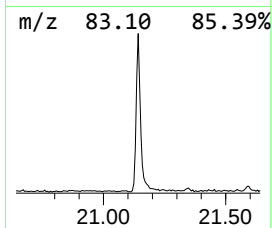
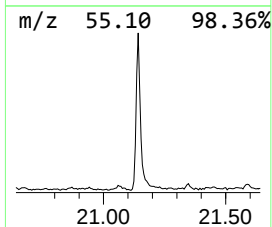
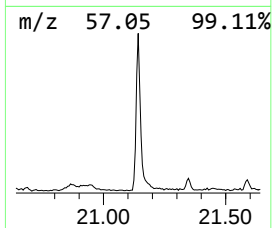
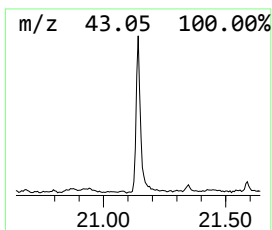
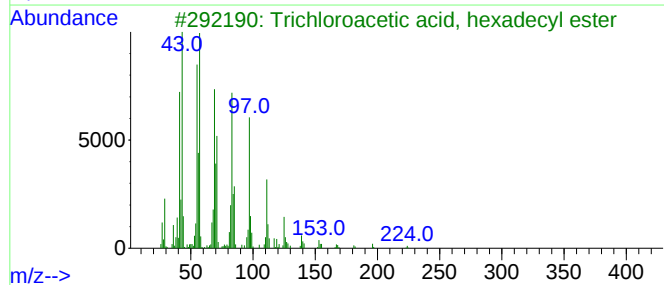
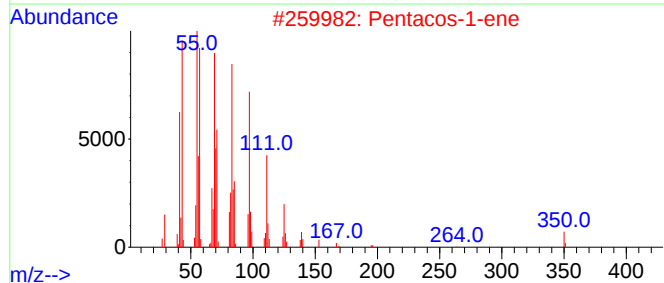
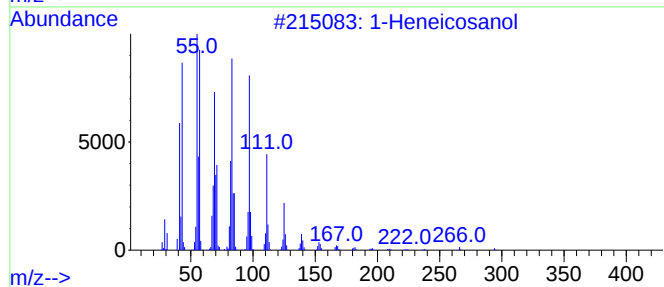
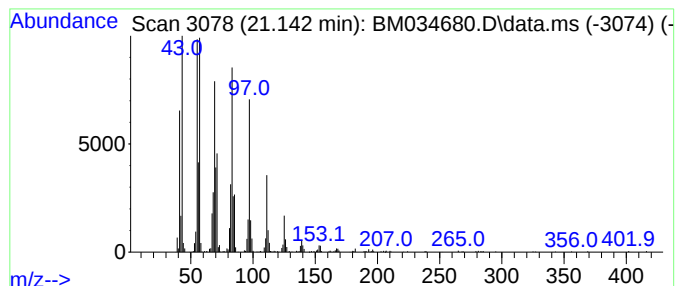
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.142	9.07 ng/ul	617127	Chrysene-d12	21.430

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	94
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			1-Tricosene	322	C23H46	018835-32-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034680.D
Acq On : 14 Apr 2022 22:50
Operator : CG/JU
Sample : N2316-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.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
Butane, 1-bromo-	4.672	3.6	ng/ul	135973	1	7.866	756087	20.0
Cyclobutane, 1,...	8.160	3.8	ng/ul	143479	1	7.866	756087	20.0
unknown-01	8.484	8.2	ng/ul	310190	1	7.866	756087	20.0
unknown-02	10.895	3.7	ng/ul	193625	2	10.672	1041520	20.0
unknown-03	11.219	2.4	ng/ul	126562	2	10.672	1041520	20.0
p-Dioxane-2,5-d...	11.630	10.2	ng/ul	530272	2	10.672	1041520	20.0
unknown-04	12.101	5.0	ng/ul	262411	2	10.672	1041520	20.0
unknown-05	12.183	4.9	ng/ul	256944	2	10.672	1041520	20.0
unknown-06	12.383	9.7	ng/ul	505147	2	10.672	1041520	20.0
n-Hexadecanoic ...	18.154	3.4	ng/ul	250202	4	17.254	1473680	20.0
1-Heneicosanol	21.142	9.1	ng/ul	617127	5	21.430	1360590	20.0