

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
 Data File : BM030476.D
 Acq On : 16 Jun 2021 21:10
 Operator : CG/JU
 Sample : M2596-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG5Q1

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-BM060621.M
 Title : SVOA CALIBRATION

Signal : TIC: BM030476.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.322	34	40	67	rBV	3034489	5833398	23.21%	2.992%
2	3.816	120	124	131	rVB	97727	141673	0.56%	0.073%
3	3.969	132	150	153	rBV3	235323	816544	3.25%	0.419%
4	5.481	401	407	415	rBV	75306	162164	0.65%	0.083%
5	6.998	654	665	678	rBV	425138	784615	3.12%	0.402%
6	7.157	684	692	699	rBV	354948	555935	2.21%	0.285%
7	7.363	721	727	737	rVB	463165	771993	3.07%	0.396%
8	7.822	796	805	813	rBV	586684	984094	3.92%	0.505%
9	8.528	920	925	932	rVV	493947	818068	3.26%	0.420%
10	8.922	983	992	996	rBV6	65313	168006	0.67%	0.086%
11	8.981	996	1002	1008	rVV	414809	727971	2.90%	0.373%
12	9.045	1008	1013	1017	rVV	118330	193189	0.77%	0.099%
13	9.463	1078	1084	1089	rVV2	82569	162902	0.65%	0.084%
14	9.539	1089	1097	1101	rVV5	68833	168533	0.67%	0.086%
15	9.698	1116	1124	1129	rBV	395786	685127	2.73%	0.351%
16	9.798	1136	1141	1148	rBV6	102382	184801	0.74%	0.095%
17	9.869	1148	1153	1155	rBV2	106511	172801	0.69%	0.089%
18	9.963	1161	1169	1174	rVB3	140608	236236	0.94%	0.121%
19	10.039	1174	1182	1194	rBV3	182089	463141	1.84%	0.238%
20	10.145	1195	1200	1204	rBV	160275	234906	0.93%	0.120%
21	10.239	1211	1216	1224	rVB	528495	941883	3.75%	0.483%
22	10.398	1238	1243	1248	rBV4	100030	181687	0.72%	0.093%
23	10.463	1248	1254	1259	rBV	745171	1164689	4.63%	0.597%
24	10.610	1269	1279	1282	rBV	808769	1827455	7.27%	0.937%
25	10.657	1282	1287	1295	rVB	5518076	8803784	35.03%	4.516%
26	10.734	1296	1300	1302	rBV3	102234	163431	0.65%	0.084%
27	10.775	1302	1307	1312	rVB	652212	1038498	4.13%	0.533%
28	10.833	1312	1317	1322	rBV5	137173	247267	0.98%	0.127%
29	11.022	1339	1349	1355	rBV8	131413	404470	1.61%	0.207%
30	11.098	1356	1362	1368	rVB3	199568	413175	1.64%	0.212%
31	11.275	1388	1392	1393	rBV2	149252	193189	0.77%	0.099%
32	11.310	1394	1398	1405	rVV6	322129	864686	3.44%	0.444%
33	11.375	1405	1409	1416	rVV4	233618	507843	2.02%	0.261%
34	11.445	1416	1421	1428	rVV3	333456	602916	2.40%	0.309%
35	11.528	1430	1435	1442	rVB7	225424	501949	2.00%	0.257%
36	11.633	1448	1453	1458	rBV	1024703	1760418	7.00%	0.903%

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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-BM060621.M
 Title : SVOA CALIBRATION

37	11.804	1477	1482	1489	rVB4	264440	507862	2.02%	0.261%
38	11.898	1490	1498	1502	rBV4	246632	513416	2.04%	0.263%
39	12.163	1541	1543	1547	rVB3	190267	245716	0.98%	0.126%
40	12.257	1552	1559	1566	rVB4	399554	1077206	4.29%	0.553%
41	12.339	1568	1573	1580	rBV7	392038	755010	3.00%	0.387%
42	12.486	1593	1598	1600	rBV4	239905	416841	1.66%	0.214%
43	12.522	1601	1604	1609	rVB3	381798	575300	2.29%	0.295%
44	12.669	1623	1629	1632	rBV3	359699	597062	2.38%	0.306%
45	12.727	1632	1639	1644	rBV6	313839	784944	3.12%	0.403%
46	12.775	1644	1647	1650	rBV2	232736	297662	1.18%	0.153%
47	12.939	1671	1675	1676	rBV3	223999	299535	1.19%	0.154%
48	12.986	1679	1683	1689	rVB	1542861	2346840	9.34%	1.204%
49	13.086	1695	1700	1707	rVB3	1022742	1782618	7.09%	0.914%
50	13.233	1721	1725	1728	rBV3	492056	829501	3.30%	0.425%
51	13.410	1745	1755	1758	rBV5	556742	1560610	6.21%	0.801%
52	13.510	1770	1772	1775	rVB3	199628	191942	0.76%	0.098%
53	13.610	1783	1789	1796	rBV	615260	1610384	6.41%	0.826%
54	13.774	1812	1817	1821	rBV	600254	929513	3.70%	0.477%
55	13.822	1822	1825	1828	rVV2	347288	488051	1.94%	0.250%
56	13.863	1828	1832	1836	rVV2	809736	1458965	5.81%	0.748%
57	13.904	1836	1839	1841	rVV	1317631	1637483	6.52%	0.840%
58	13.933	1841	1844	1848	rVB	2218381	2785801	11.08%	1.429%
59	13.974	1848	1851	1854	rBV4	280561	342517	1.36%	0.176%
60	14.098	1869	1872	1876	rBV5	312868	526947	2.10%	0.270%
61	14.145	1876	1880	1884	rVV	1126915	1555331	6.19%	0.798%
62	14.286	1900	1904	1910	rVB3	646763	1136968	4.52%	0.583%
63	14.351	1910	1915	1920	rBV4	704668	1258099	5.01%	0.645%
64	14.421	1923	1927	1928	rVV3	709309	890048	3.54%	0.457%
65	14.451	1929	1932	1938	rVB	1441933	2076796	8.26%	1.065%
66	14.657	1963	1967	1976	rVB2	825584	2116014	8.42%	1.085%
67	14.845	1995	1999	2007	rVB3	1260556	2529407	10.06%	1.297%
68	14.921	2010	2012	2016	rVB	602944	730125	2.91%	0.375%
69	14.969	2016	2020	2030	rVB4	835035	1701919	6.77%	0.873%
70	15.069	2033	2037	2041	rBV	694044	1214587	4.83%	0.623%
71	15.298	2073	2076	2080	rVB	678548	843329	3.36%	0.433%
72	15.439	2096	2100	2105	rVB	1309785	1816511	7.23%	0.932%
73	15.486	2105	2108	2112	rBV2	391484	499422	1.99%	0.256%
74	15.704	2141	2145	2152	rVB	2786841	4166098	16.58%	2.137%
75	15.863	2168	2172	2175	rBV5	425859	706998	2.81%	0.363%
76	15.916	2178	2181	2184	rVB2	510280	650910	2.59%	0.334%

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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

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
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77	16.180	2219	2226	2238	rVB	5047305	9646649	38.38%	4.948%
78	16.598	2293	2297	2303	rVB4	761332	985111	3.92%	0.505%
79	16.839	2335	2338	2340	rBV2	1005116	1314327	5.23%	0.674%
80	16.868	2340	2343	2347	rVB2	1466828	1818116	7.23%	0.933%
81	16.945	2353	2356	2360	rVB	796241	889560	3.54%	0.456%
82	16.998	2361	2365	2370	rBV	3129667	4421383	17.59%	2.268%
83	17.186	2393	2397	2401	rBV	1724715	2490079	9.91%	1.277%
84	17.286	2410	2414	2419	rVB	1763986	2353150	9.36%	1.207%
85	17.357	2423	2426	2428	rBV2	506371	705924	2.81%	0.362%
86	17.415	2434	2436	2441	rVB3	876457	1068793	4.25%	0.548%
87	17.604	2464	2468	2477	rVV2	1080725	1916069	7.62%	0.983%
88	17.680	2478	2481	2487	rVB2	1082692	1375488	5.47%	0.706%
89	17.851	2507	2510	2515	rVB	2103495	2693425	10.72%	1.382%
90	18.104	2546	2553	2561	rVB2	12313102	20987543	83.51%	10.766%
91	18.374	2596	2599	2604	rVB	712751	839977	3.34%	0.431%
92	18.974	2699	2701	2704	rBV	558545	685691	2.73%	0.352%
93	19.156	2729	2732	2741	rVB2	2305411	2987425	11.89%	1.532%
94	19.239	2743	2746	2750	rVB	782953	989136	3.94%	0.507%
95	19.368	2762	2768	2784	rBV	11371765	18818380	74.88%	9.653%
96	19.574	2799	2803	2806	rBV3	2038425	3012229	11.99%	1.545%
97	21.050	3048	3054	3060	rBV	18279487	25131890	100.00%	12.892%
98	23.327	3436	3441	3453	rVB	2932087	5345497	21.27%	2.742%
99	23.480	3464	3467	3473	rVB	1061955	1597683	6.36%	0.820%
100	23.627	3488	3492	3498	rVB2	1444799	2536943	10.09%	1.301%

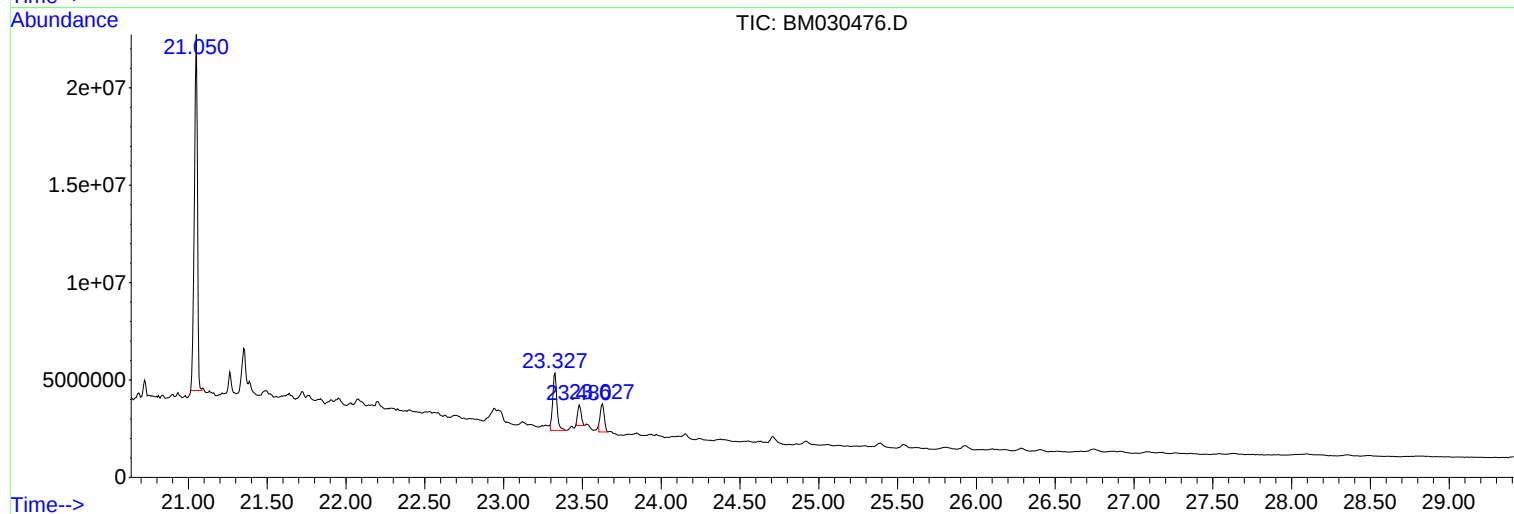
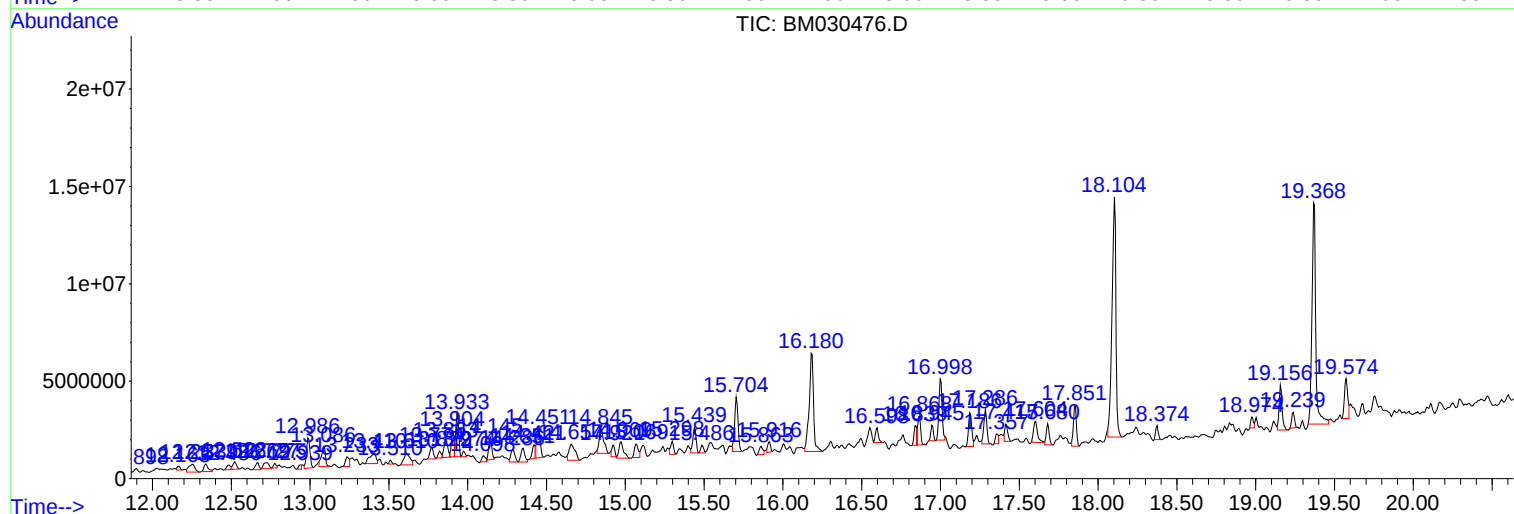
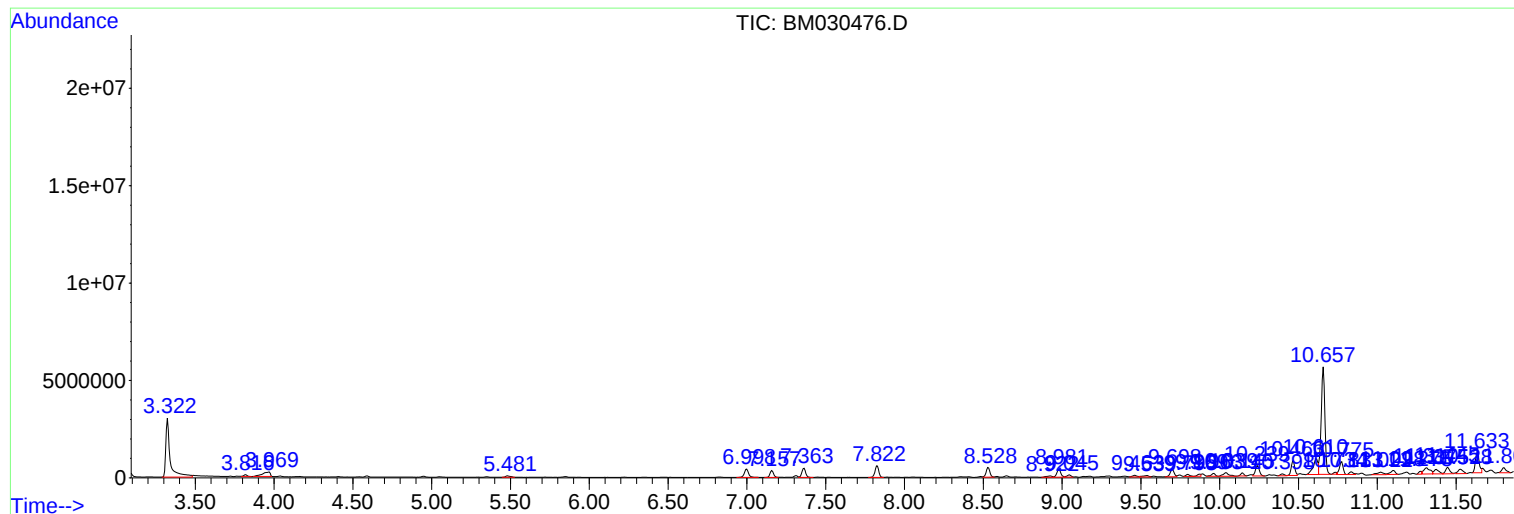
Sum of corrected areas: 194948193

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Instrument :
BNA_M
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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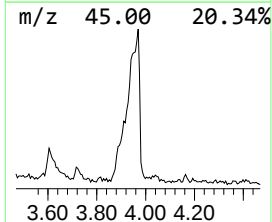
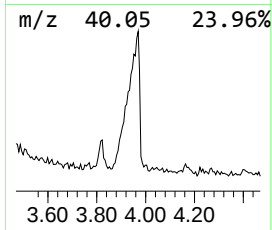
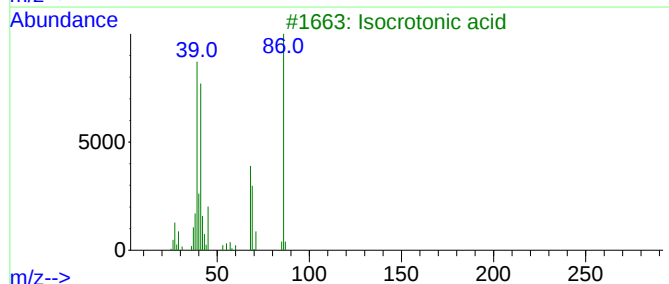
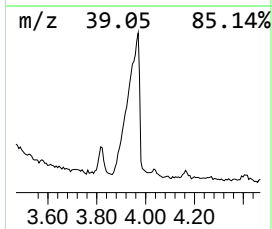
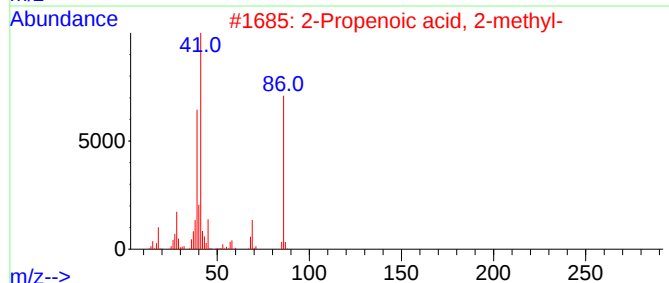
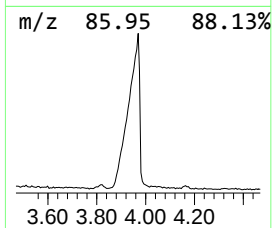
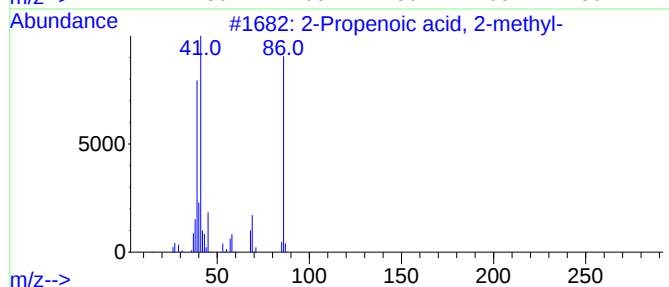
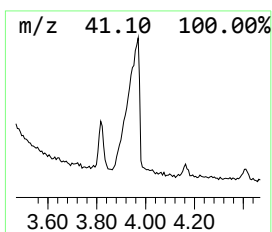
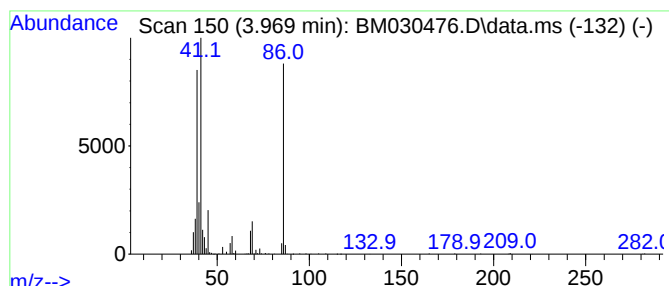
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Propenoic acid, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.970	16.59 ng/ul	816544	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, 2-methyl-		86	C4H6O2		000079-41-4	96
2	2-Propenoic acid, 2-methyl-		86	C4H6O2		000079-41-4	94
3	Isocrotonic acid		86	C4H6O2		000503-64-0	91
4	2-Propenoic acid, 2-methyl-		86	C4H6O2		000079-41-4	91
5	2-Butenoic acid, (E)-		86	C4H6O2		000107-93-7	91



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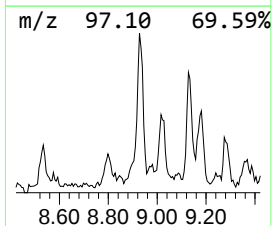
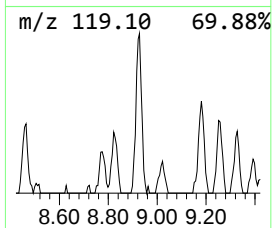
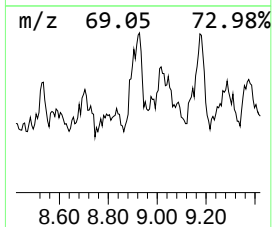
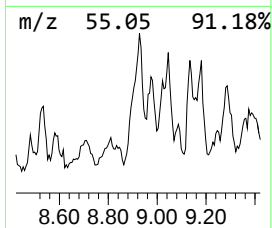
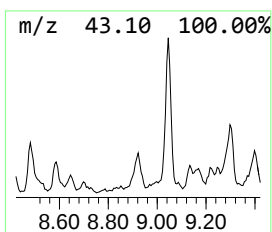
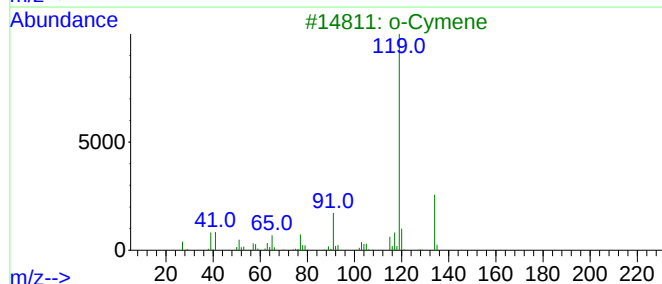
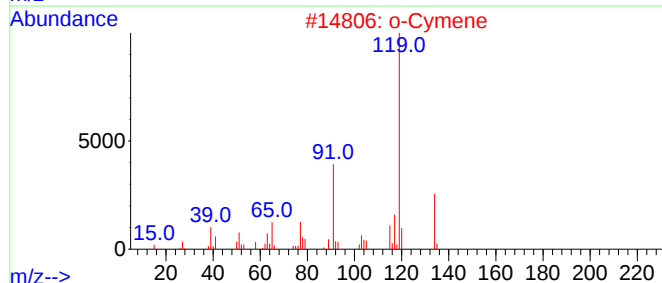
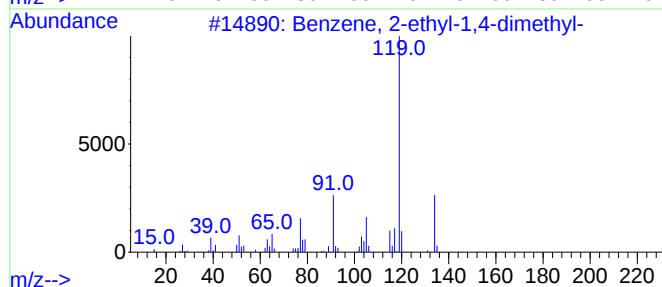
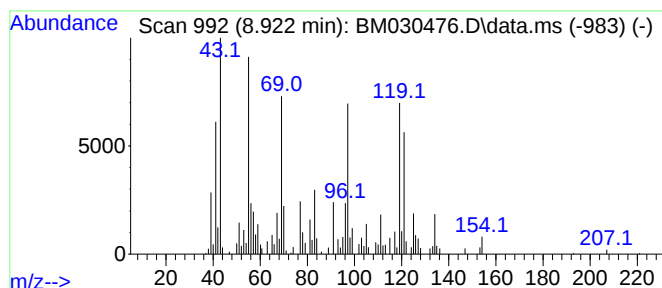
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-01 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.920	3.41 ng/ul	168006	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	42
2			o-Cymene	134	C10H14	000527-84-4	42
3			o-Cymene	134	C10H14	000527-84-4	42
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	42
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	42



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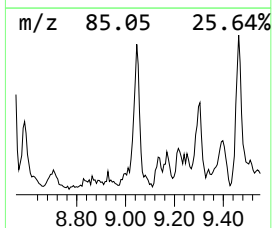
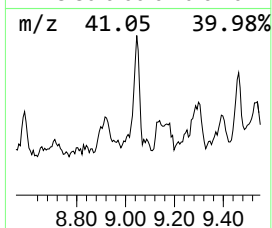
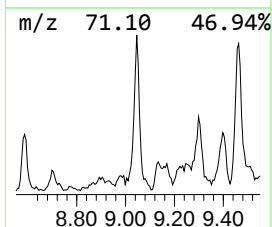
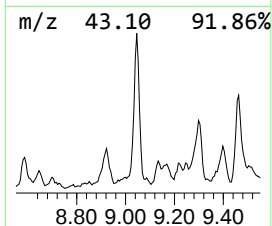
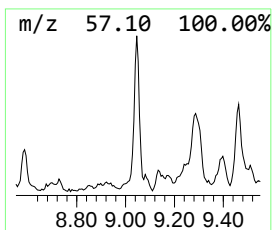
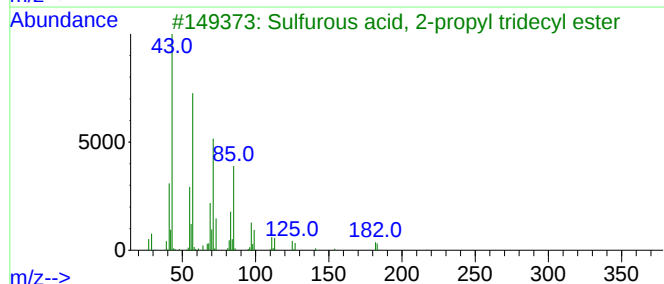
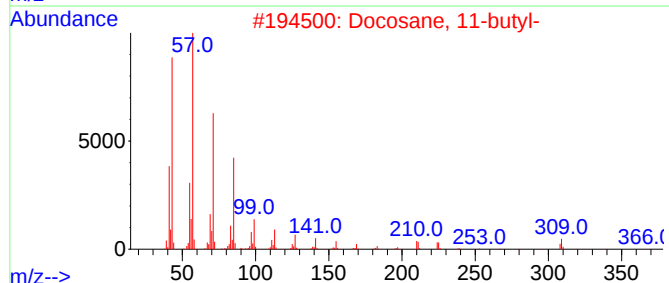
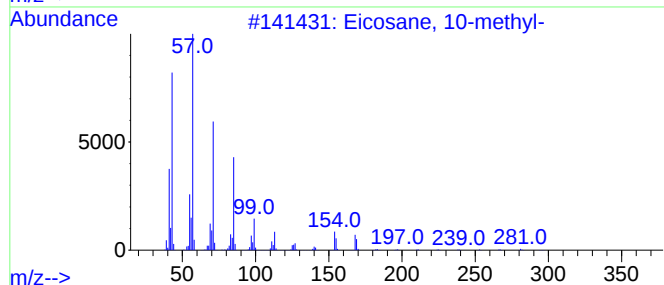
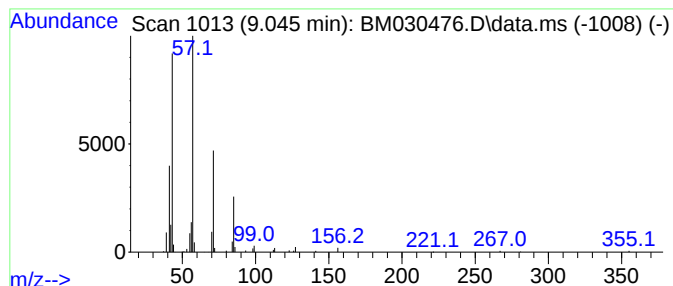
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.050	3.93 ng/ul	193189	1,4-Dichlorobenzene-d4	7.828	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	83
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	78
3	Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	78
4	Tridecane	184	C13H28	000629-50-5	78
5	Tricosane, 2-methyl-	338	C24H50	001928-30-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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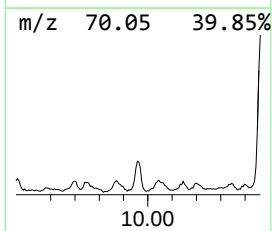
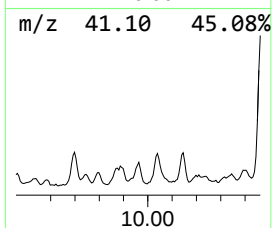
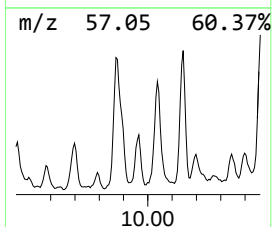
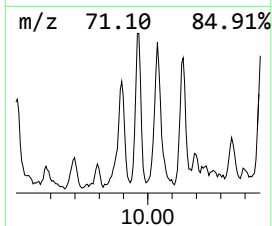
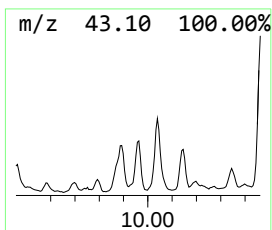
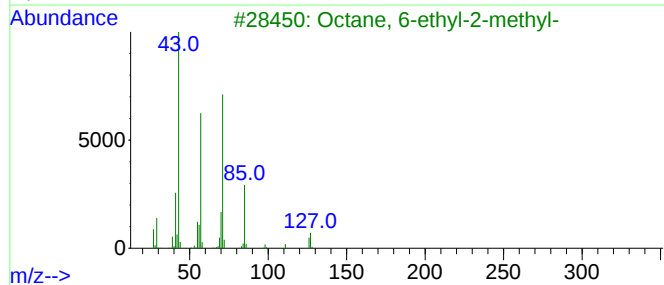
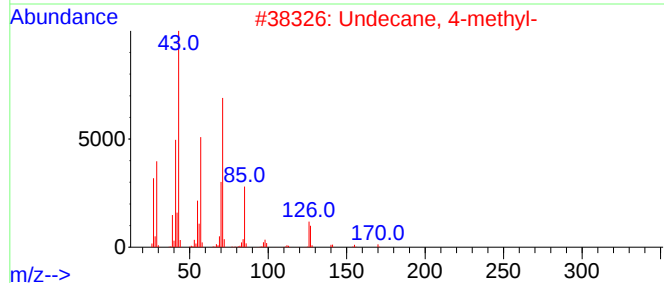
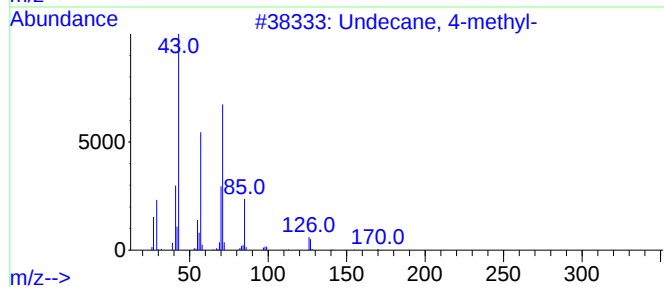
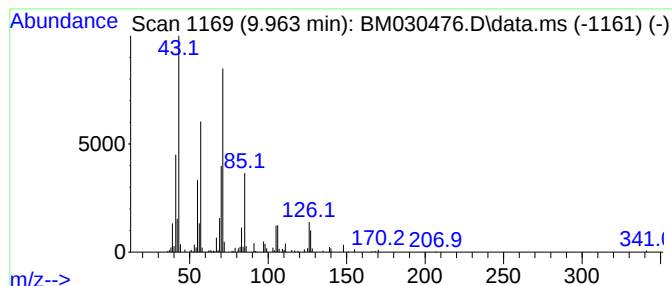
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.960	2.59 ng/ul	236236	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 4-methyl-	170	C12H26	002980-69-0	87
2		Undecane, 4-methyl-	170	C12H26	002980-69-0	76
3		Octane, 6-ethyl-2-methyl-	156	C11H24	062016-19-7	50
4		Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	47
5		n-Amyl ether	158	C10H22O	000693-65-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

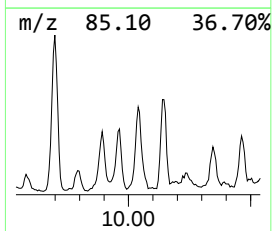
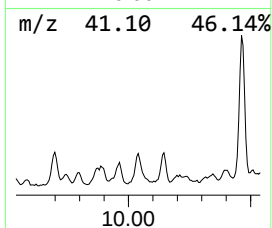
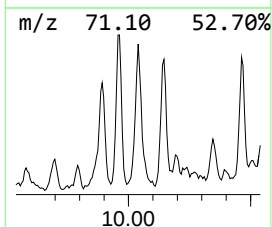
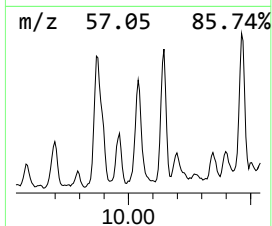
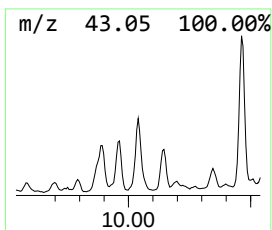
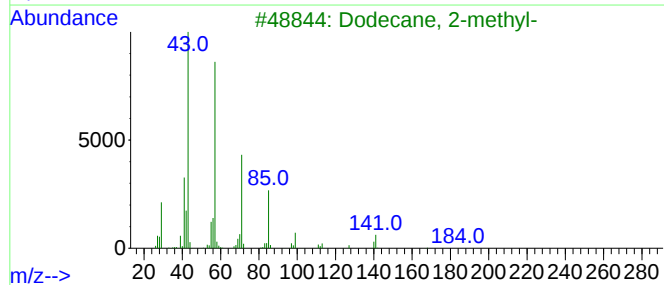
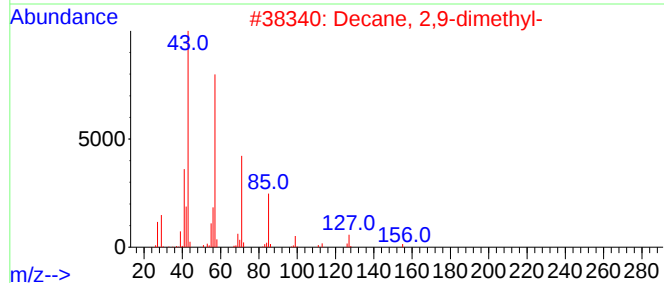
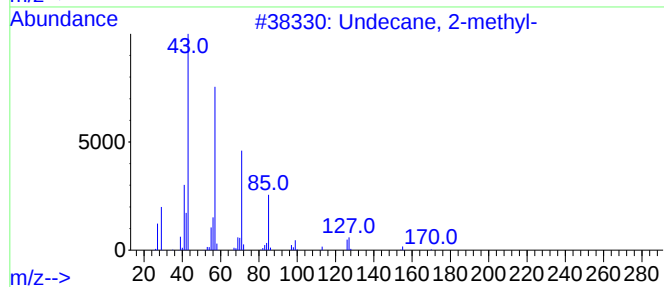
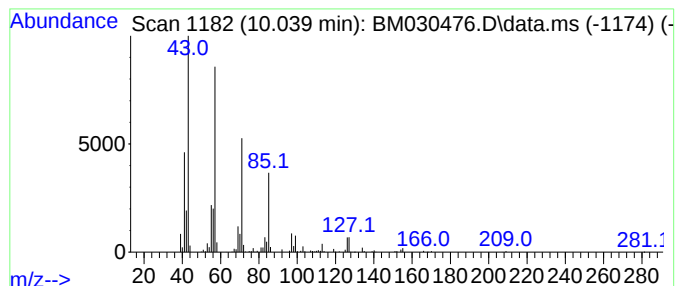
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.040	5.07 ng/ul	463141	Naphthalene-d8	10.610	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2-methyl-	170	C12H26	007045-71-8	91
2	Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	83
3	Dodecane, 2-methyl-	184	C13H28	001560-97-0	78
4	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	72
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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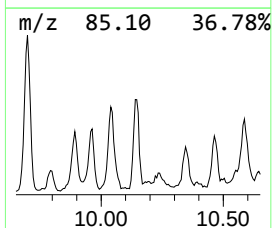
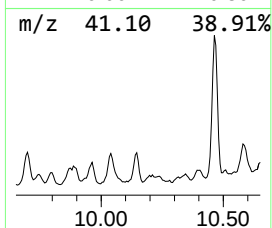
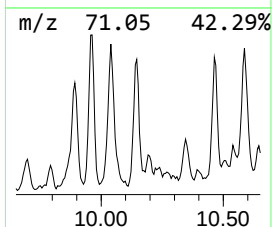
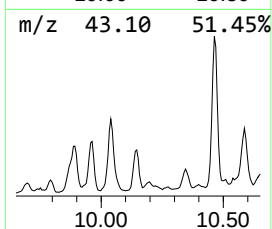
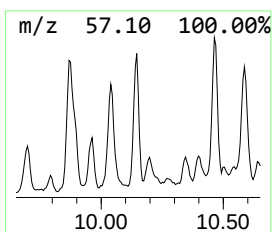
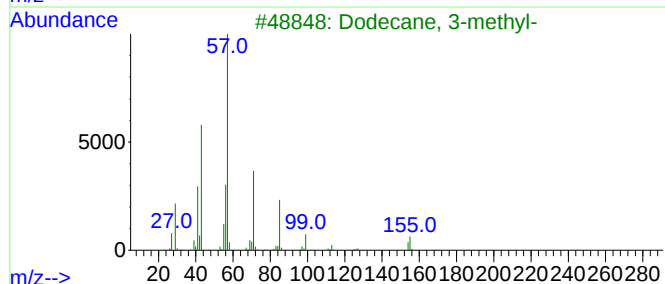
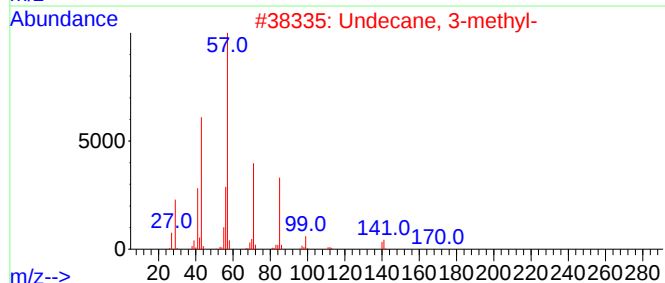
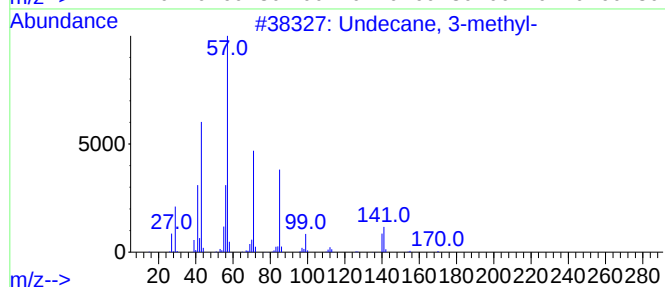
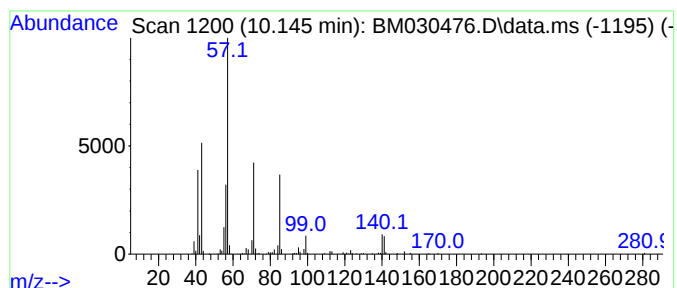
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.150	2.57 ng/ul	234906	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3-methyl-	170	C12H26	001002-43-3	91
2		Undecane, 3-methyl-	170	C12H26	001002-43-3	83
3		Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
4		Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64
5		Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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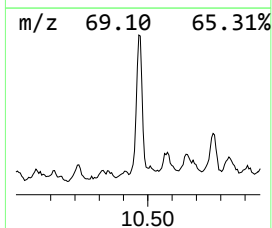
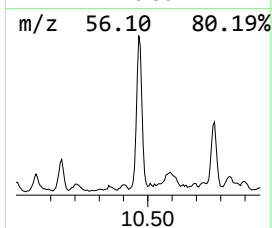
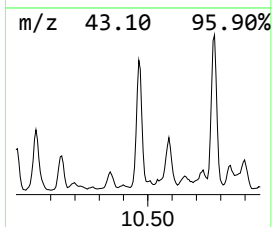
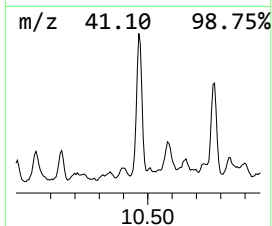
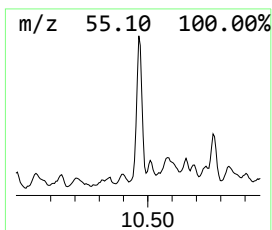
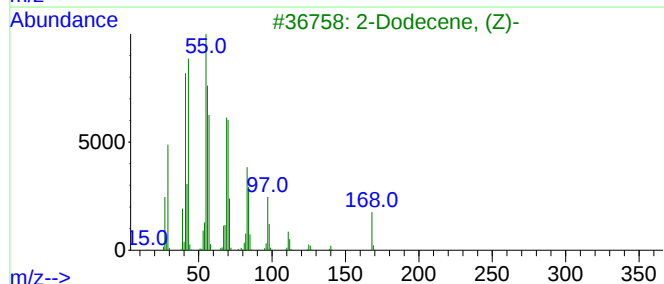
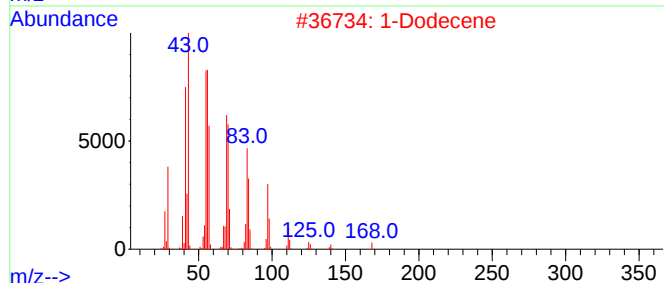
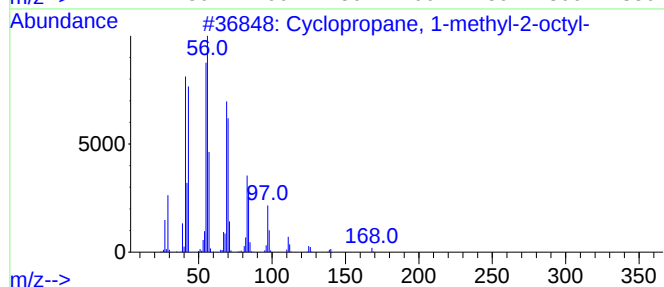
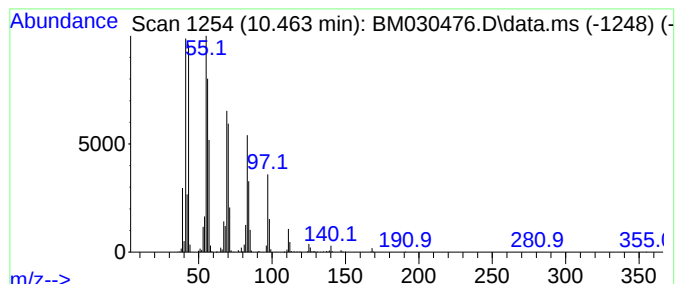
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Cyclic10.46 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.460	12.75 ng/ul	1164690	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, 1-methyl-2-octyl-	168	C12H24	037617-26-8	93
2		1-Dodecene	168	C12H24	000112-41-4	93
3		2-Dodecene, (Z)-	168	C12H24	007206-26-0	90
4		3-Dodecene, (Z)-	168	C12H24	007239-23-8	90
5		2-Dodecene, (Z)-	168	C12H24	007206-26-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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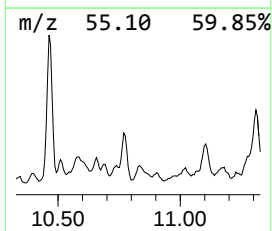
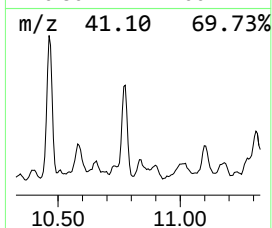
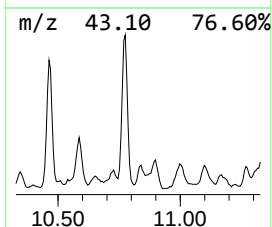
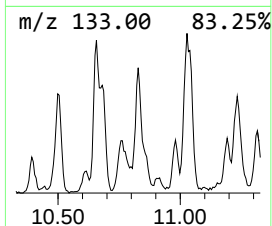
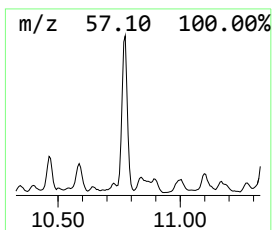
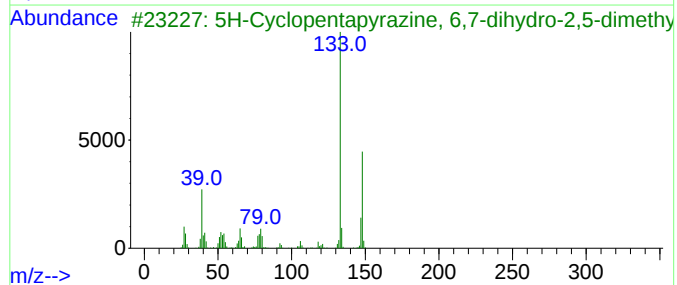
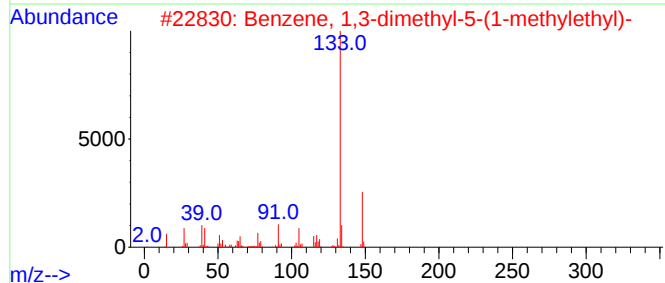
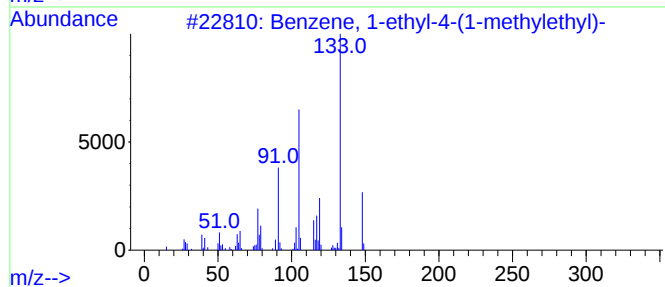
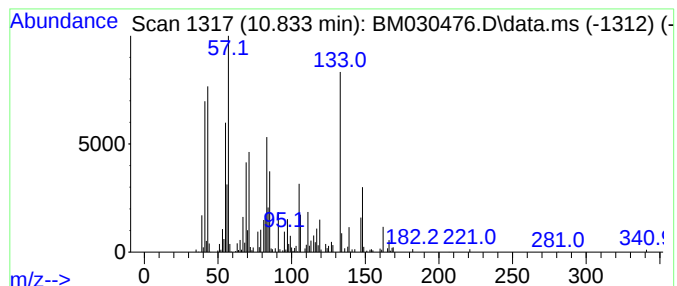
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown-02 Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.830	2.71 ng/ul	247267	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	42
2			Benzene, 1,3-dimethyl-5-(1-methylethyl)-	148	C11H16	004706-90-5	30
3			5H-Cyclopentapyrazine, 6,7-dihydro-2,5-dimethyl-	148	C9H12N2	038917-61-2	30
4			3,4-Dimethylcumene	148	C11H16	1000370-34-1	30
5			Benzene, pentamethyl-	148	C11H16	000700-12-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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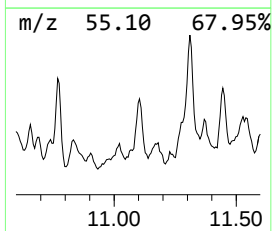
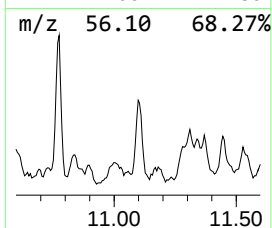
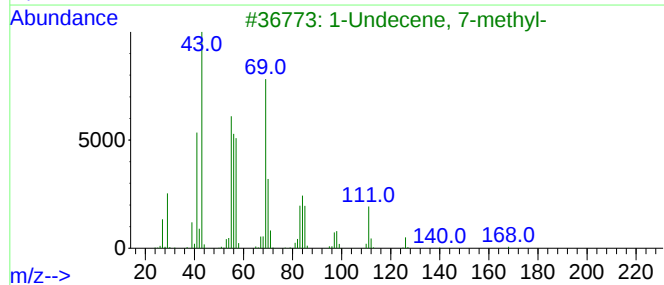
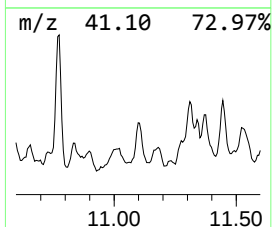
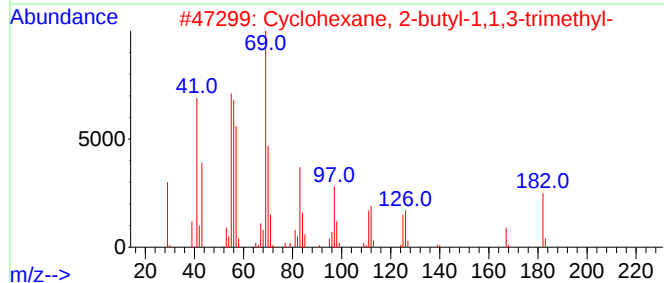
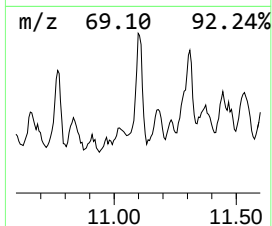
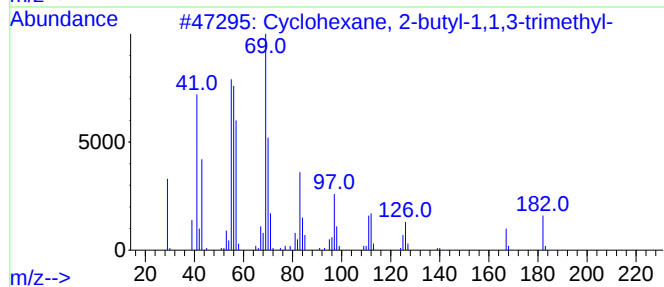
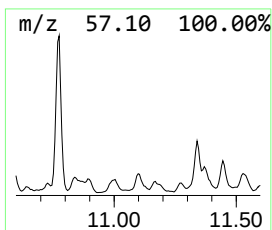
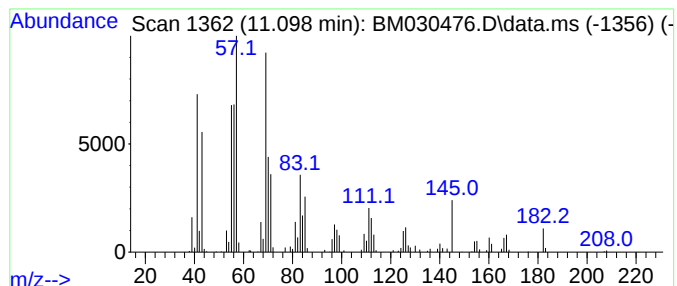
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Cyclic11.10 Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.100	4.52 ng/ul	413175	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	91
2			Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	90
3			1-Undecene, 7-methyl-	168	C12H24	074630-42-5	81
4			Cyclopentane, 1-butyl-2-propyl-	168	C12H24	062199-50-2	64
5			Cyclohexane, 1,1-dimethyl-2-propyl-	154	C11H22	081983-71-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Misc :
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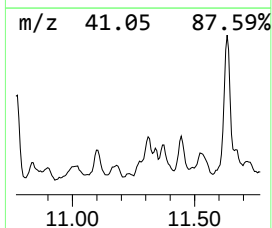
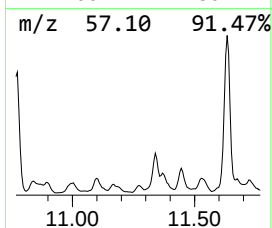
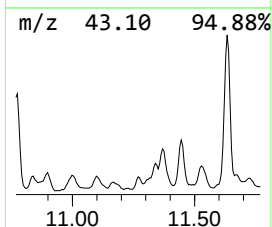
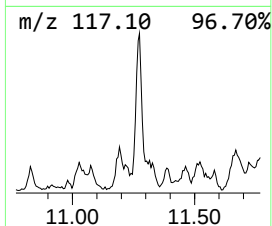
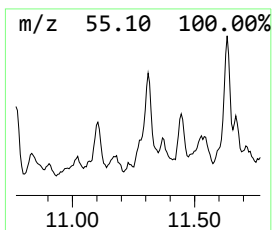
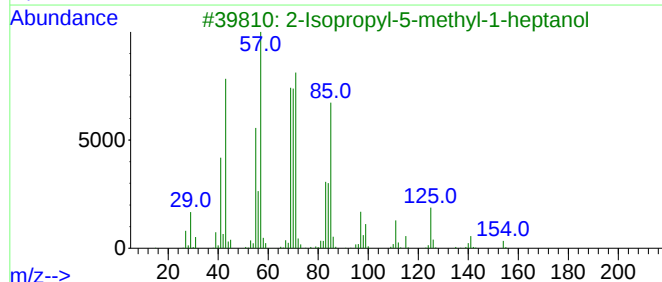
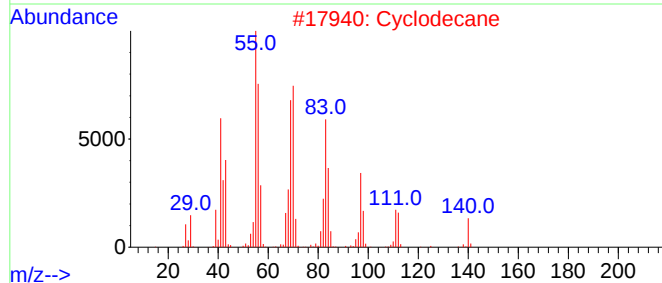
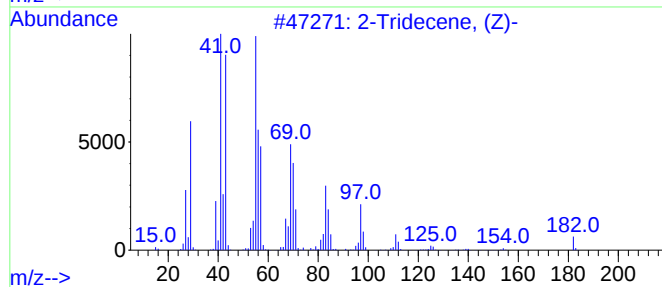
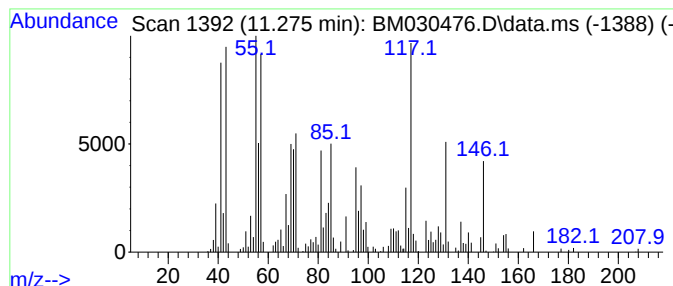
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.270	2.11 ng/ul	193189	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Tridecene, (Z)-			182	C13H26	041446-59-7	25
2	Cyclodecane			140	C10H20	000293-96-9	25
3	2-Isopropyl-5-methyl-1-heptanol			172	C11H24O	091337-07-4	25
4	1H-Indene-5-carboxylic acid, 2,3...			162	C10H10O2	1000337-61-3	25
5	1H-Indene, 1-ethyl-2,3-dihydro-			146	C11H14	004830-99-3	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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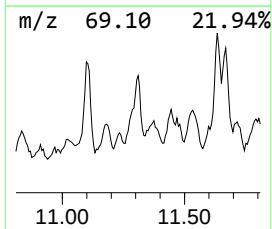
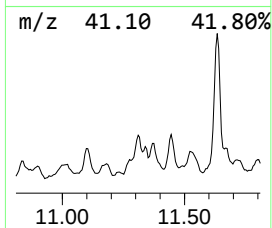
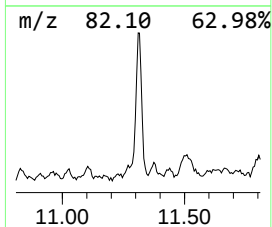
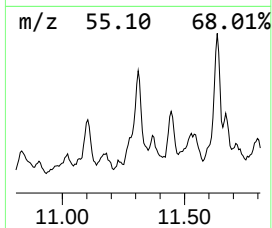
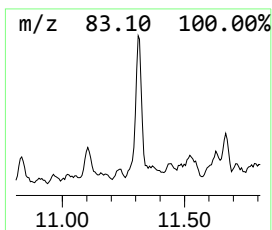
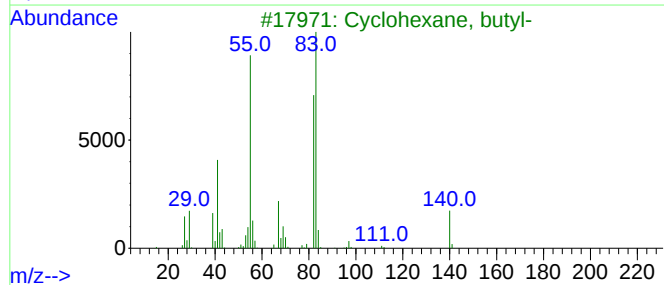
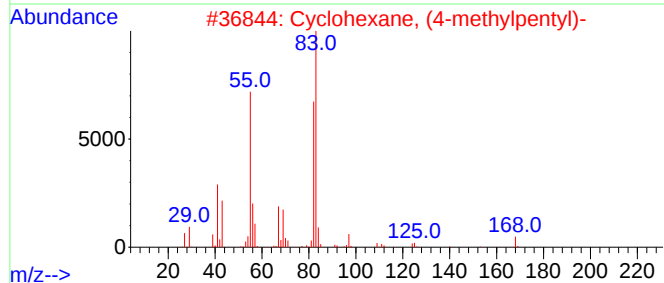
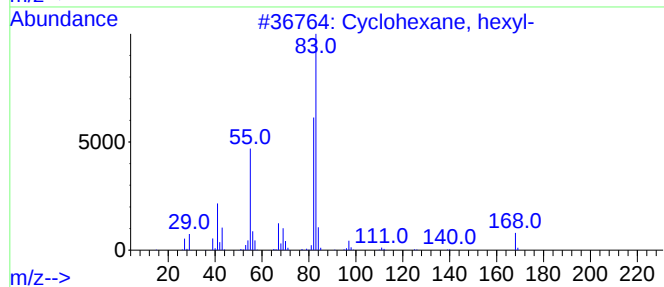
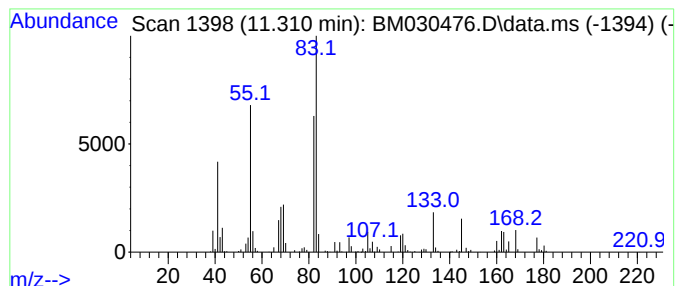
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 (DEL) Alkane: Cyclic11.31 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.310	9.46 ng/ul	864686	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, hexyl-	168	C12H24	004292-75-5	52
2			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	50
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	47
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	47
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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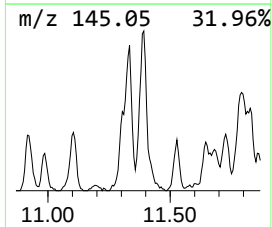
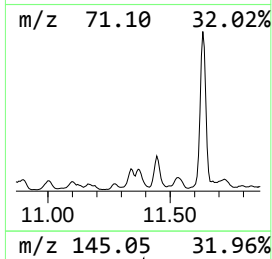
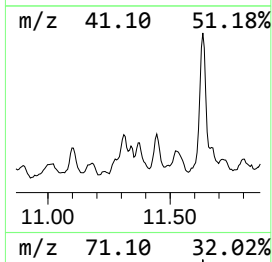
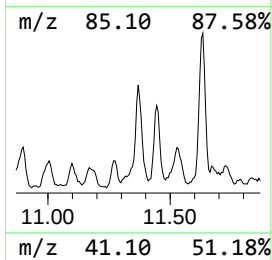
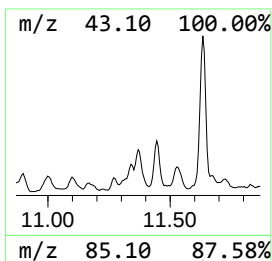
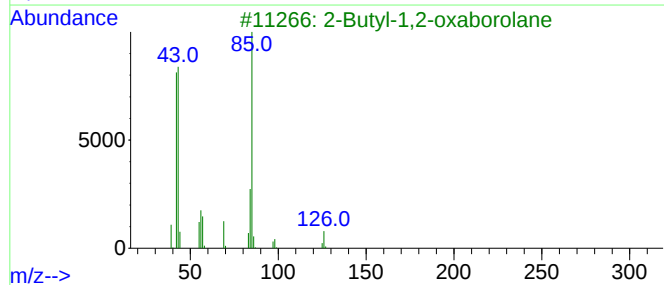
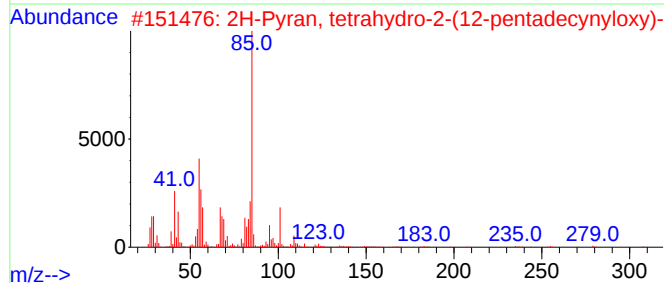
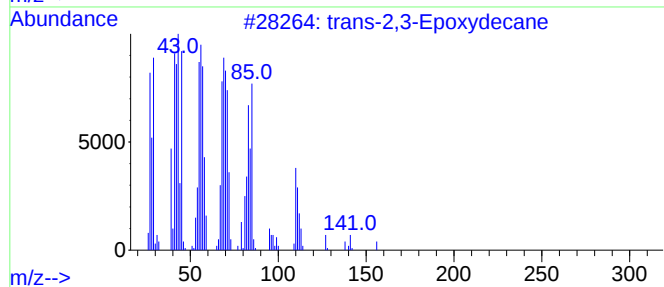
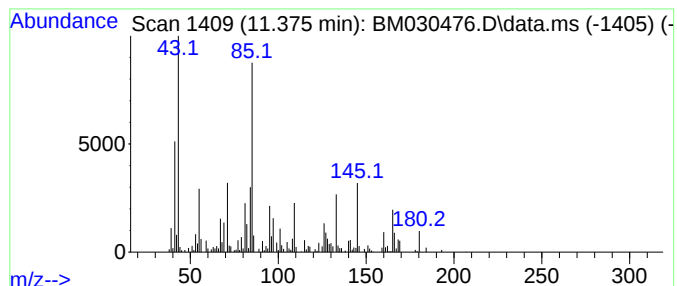
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown-03 Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.370	5.56 ng/ul	507843	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	38
2			2H-Pyran, tetrahydro-2-(12-penta...	308	C20H36O2	056666-38-7	35
3			2-Butyl-1,2-oxaborolane	126	C7H15BO	005357-13-1	30
4			2-Pentanone, 3-ethyl-3-methyl-	128	C8H16O	019780-65-5	27
5			2-Pentanone, 3-ethyl-3-methyl-	128	C8H16O	019780-65-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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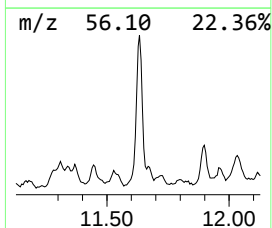
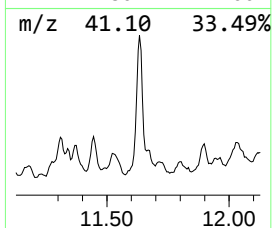
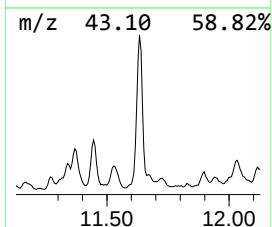
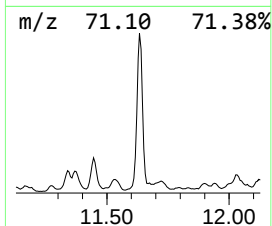
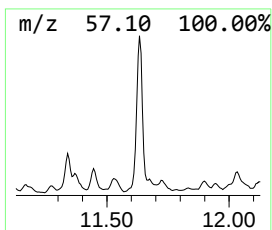
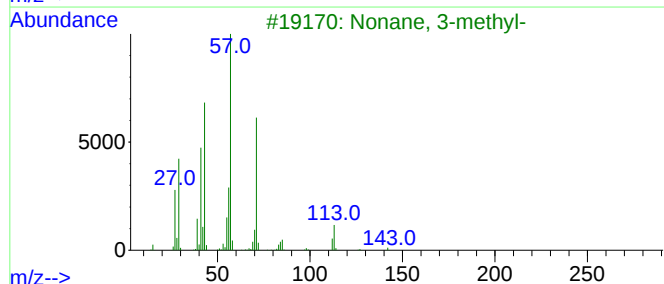
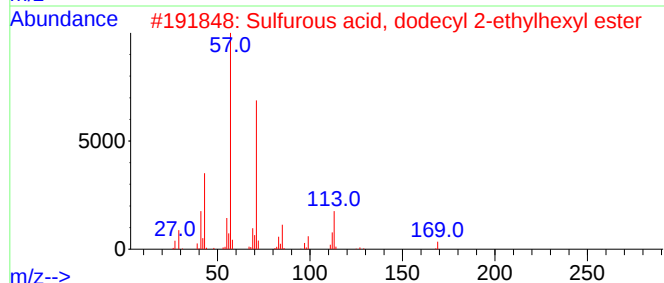
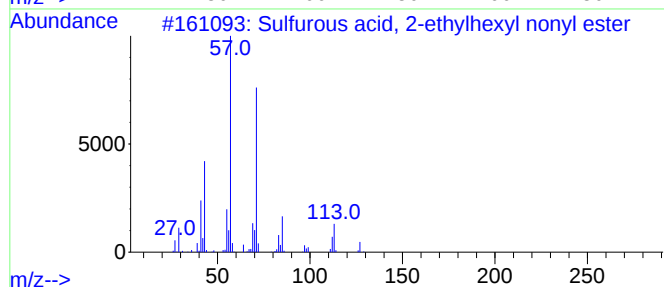
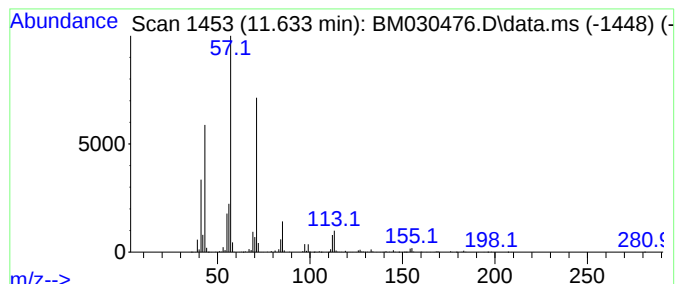
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Sulfurous acid, 2-ethylhexy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.630	19.27 ng/ul	1760420	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	72
2			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	72
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	72
4			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	72
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

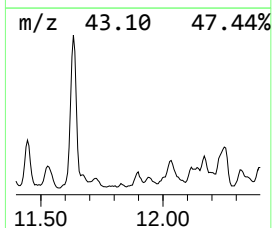
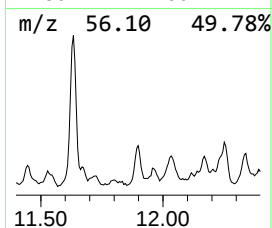
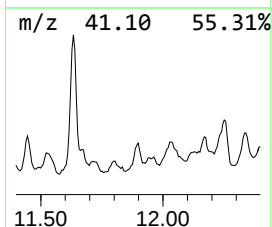
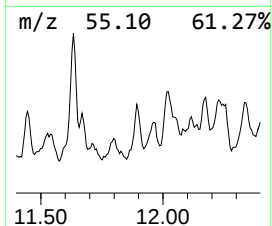
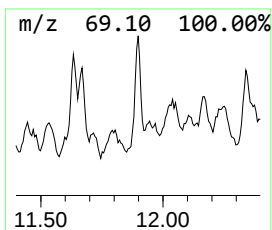
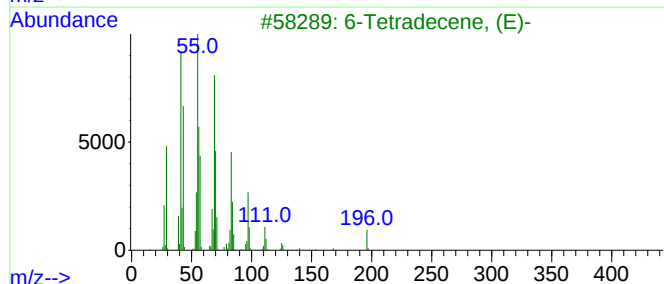
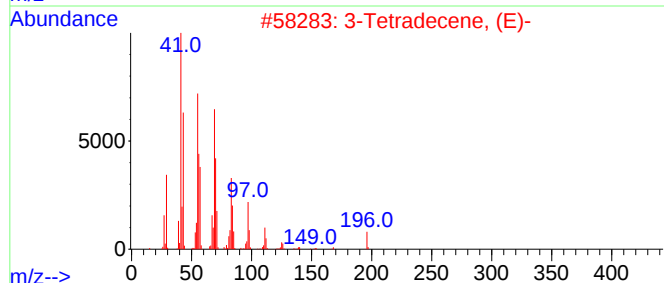
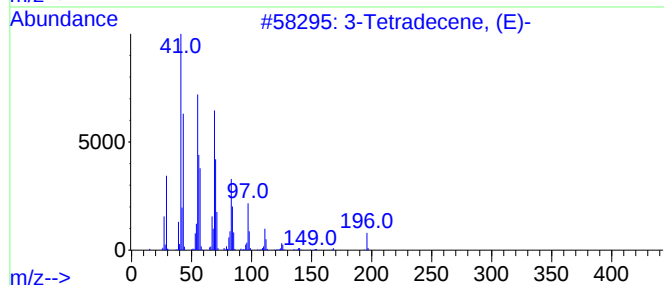
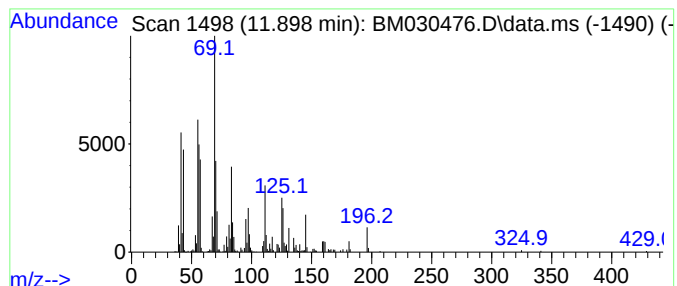
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.900	5.62 ng/ul	513416	Naphthalene-d8	10.610	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Tetradecene, (E)-	196	C14H28	041446-68-8	70
2	3-Tetradecene, (E)-	196	C14H28	041446-68-8	70
3	6-Tetradecene, (E)-	196	C14H28	041446-64-4	55
4	4-Tetradecene, (E)-	196	C14H28	041446-78-0	50
5	7-Tetradecene, (E)-	196	C14H28	041446-63-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5Q1

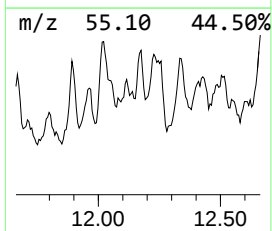
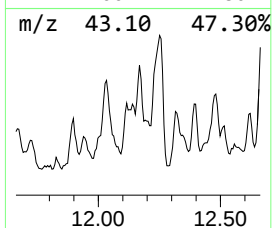
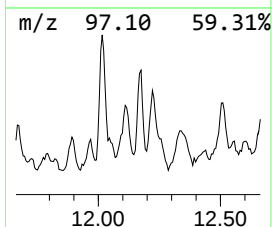
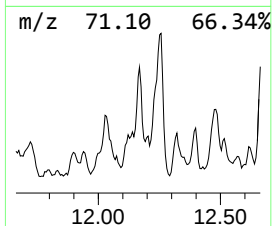
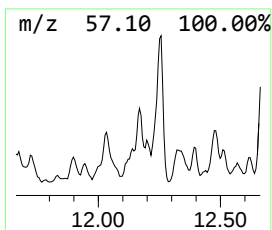
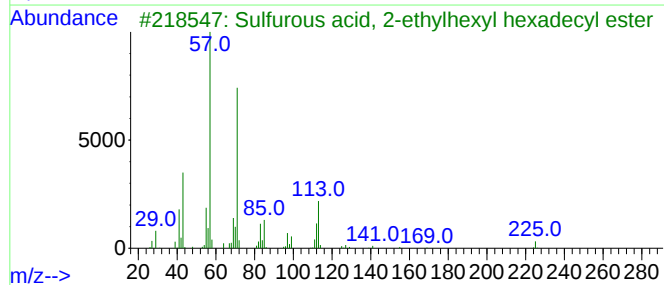
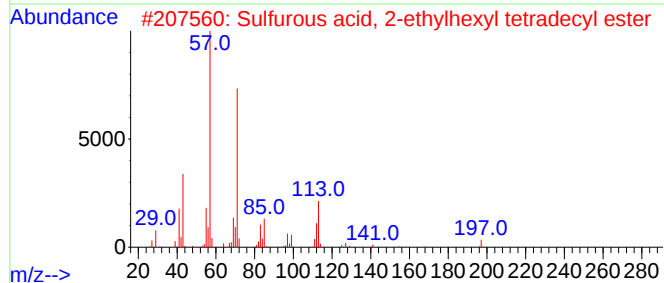
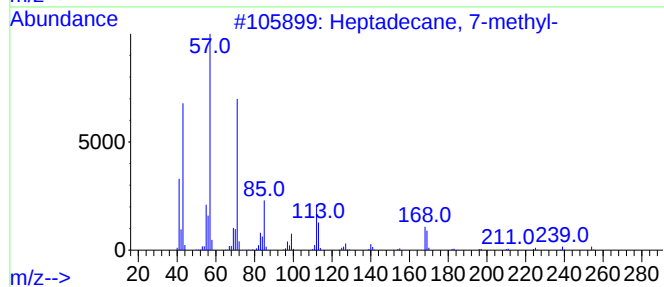
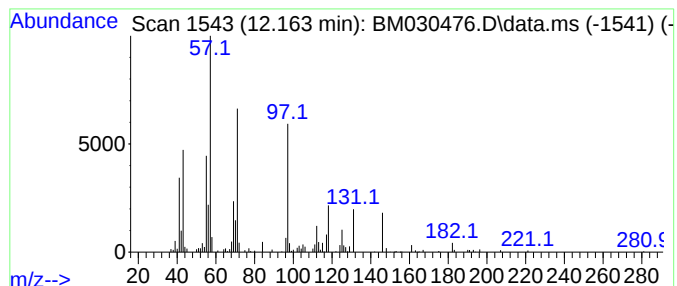
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.160	2.69 ng/ul	245716	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	35
2			Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	35
3			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	35
4			Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	27
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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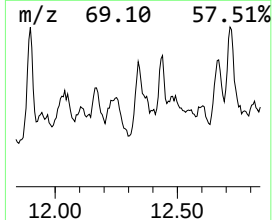
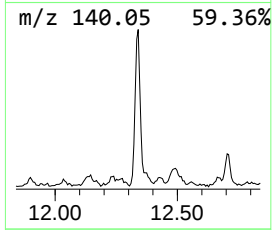
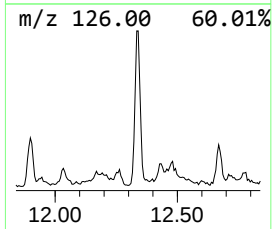
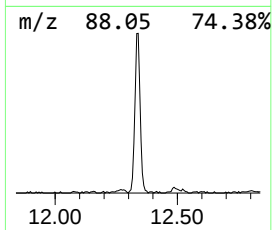
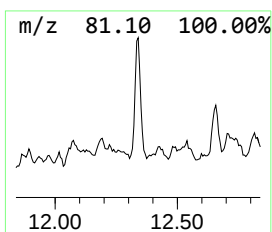
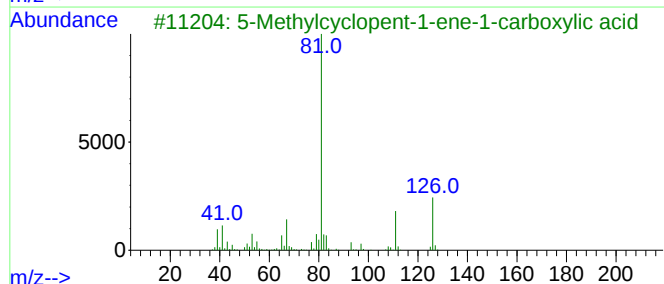
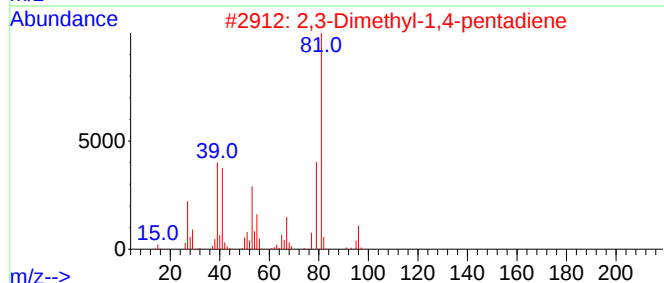
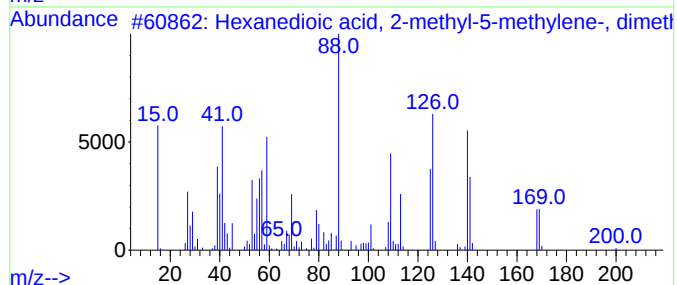
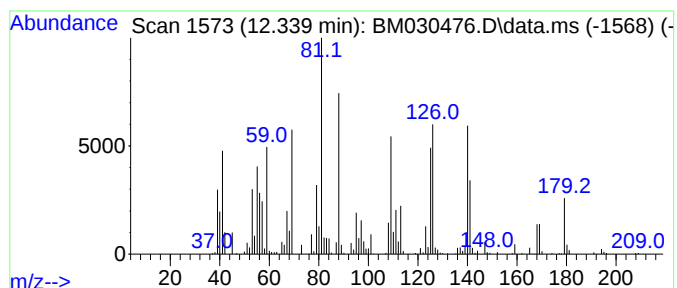
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Hexanedioic acid, 2-methyl-... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.340	8.26 ng/ul	755010	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, 2-methyl-5-met...	200	C10H16O4	004513-62-6	64
2			2,3-Dimethyl-1,4-pentadiene	96	C7H12	000758-86-1	15
3			5-Methylcyclopent-1-ene-1-carbox...	126	C7H10O2	053623-50-0	14
4			1,4-Hexadiene, 3-ethyl-	110	C8H14	002080-89-9	11
5			2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

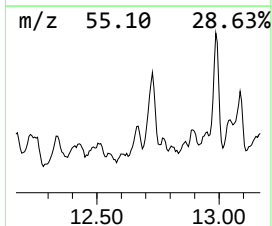
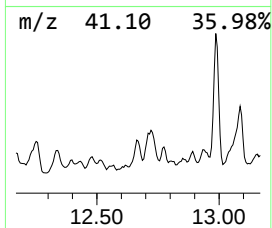
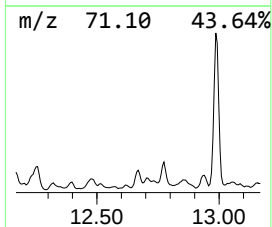
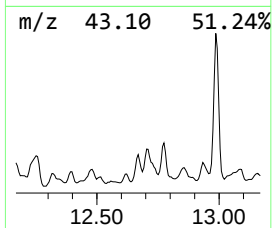
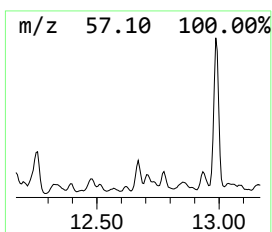
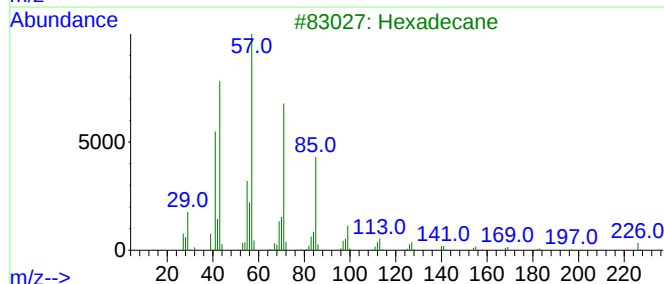
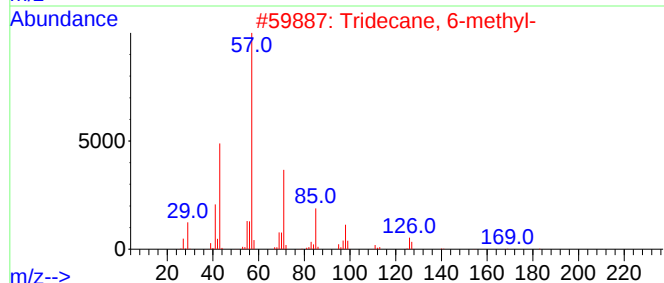
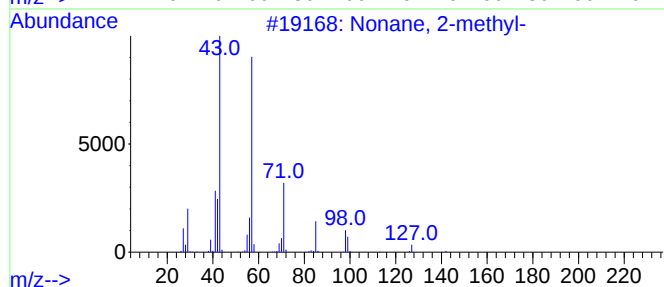
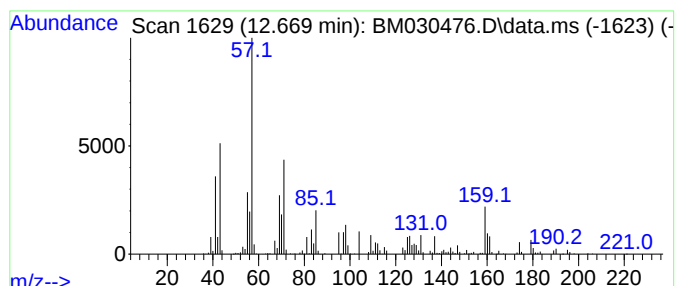
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.670	5.75 ng/ul	597062	Acenaphthene-d10		14.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-		142	C10H22	000871-83-0	52
2	Tridecane, 6-methyl-		198	C14H30	013287-21-3	49
3	Hexadecane		226	C16H34	000544-76-3	38
4	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	38
5	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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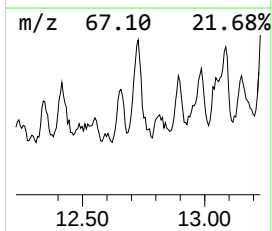
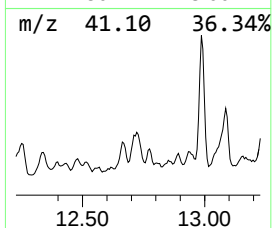
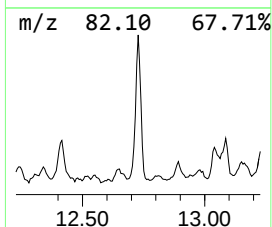
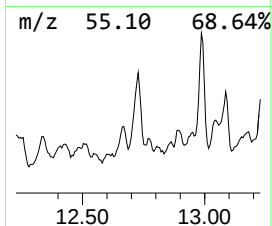
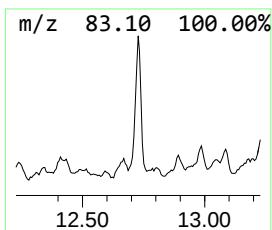
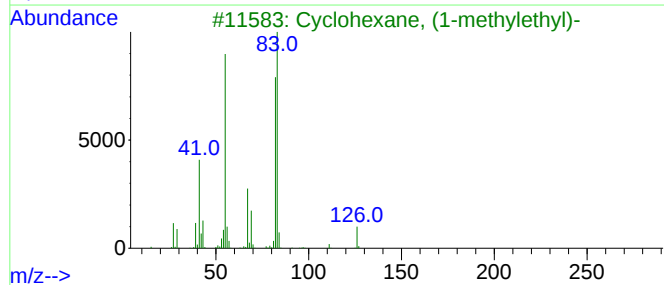
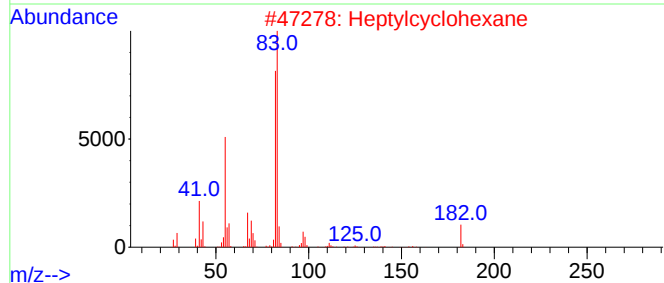
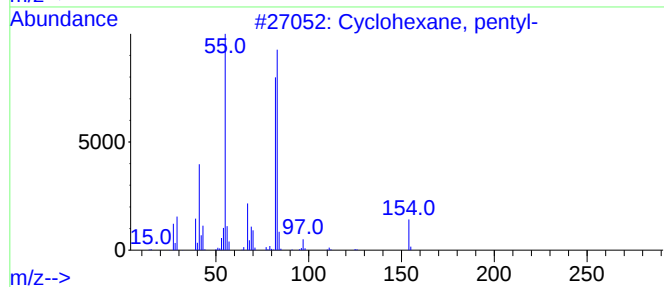
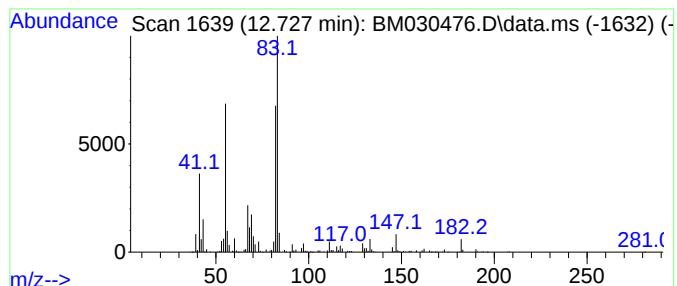
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Cyclic12.73 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.730	7.56 ng/ul	784944	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	80
2			Heptylcyclohexane	182	C13H26	005617-41-4	80
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	80
4			Cyclohexane, octyl-	196	C14H28	001795-15-9	72
5			Heptylcyclohexane	182	C13H26	005617-41-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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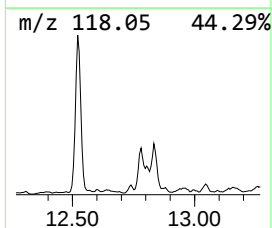
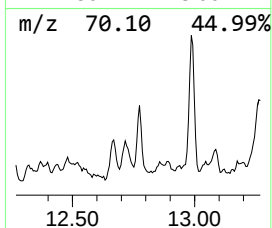
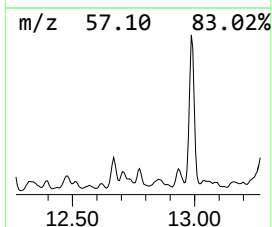
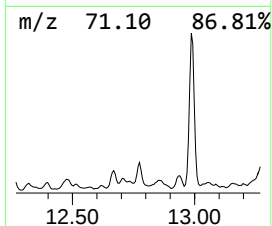
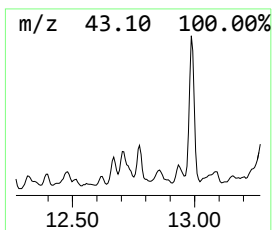
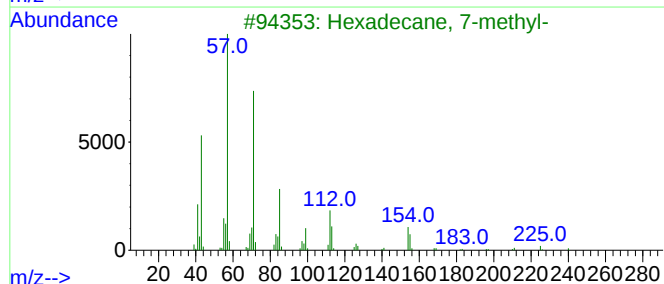
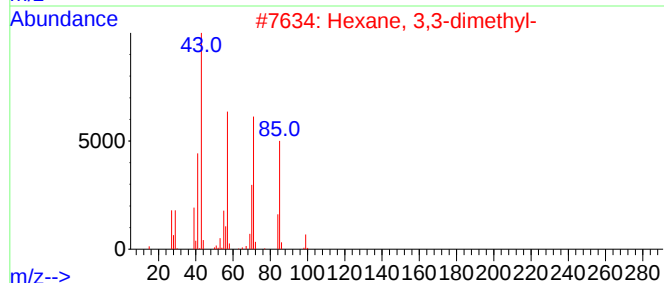
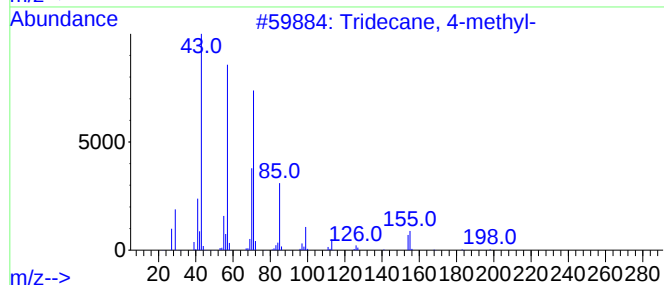
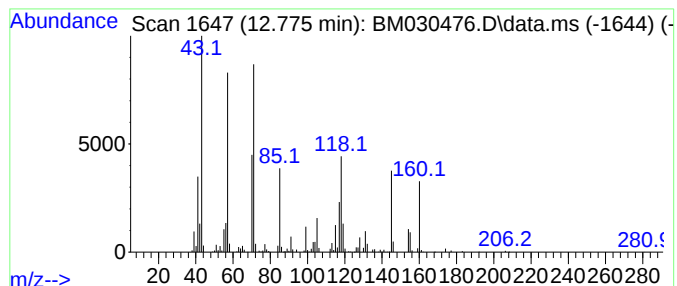
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.770	2.87 ng/ul	297662	Acenaphthene-d10	14.451

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 4-methyl-	198	C14H30	026730-12-1	49
2		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	46
3		Hexadecane, 7-methyl-	240	C17H36	026730-20-1	38
4		Sulfurous acid, decyl pentyl ester	292	C15H32O3S	1000309-14-3	35
5		Undecane, 2,8-dimethyl-	184	C13H28	017301-25-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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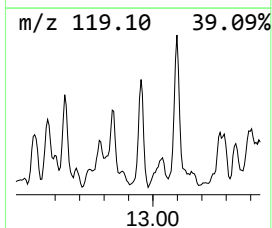
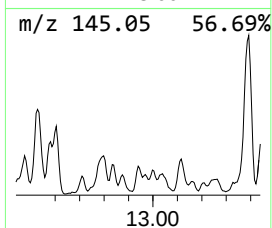
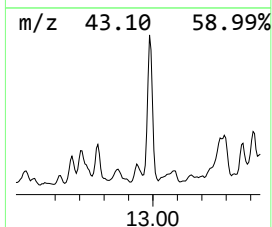
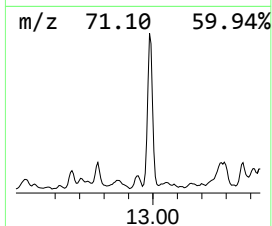
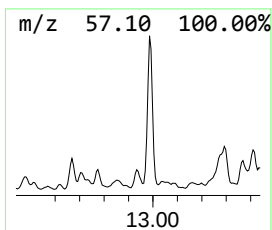
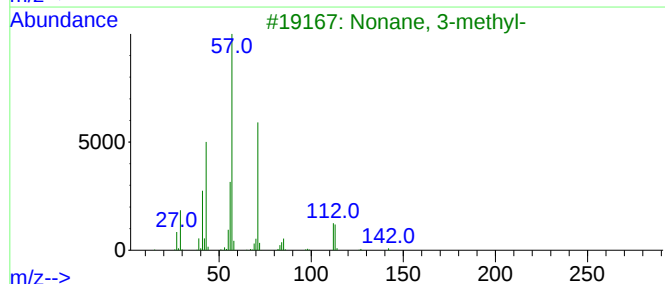
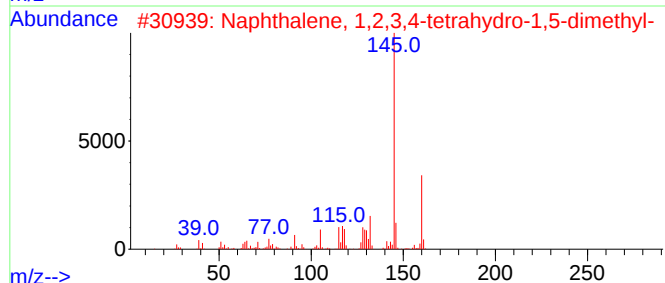
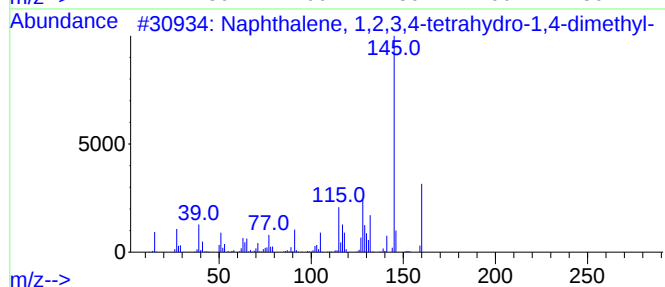
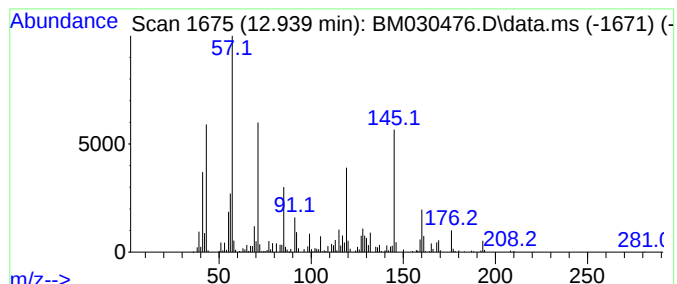
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 unknown-04 Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.940	2.88 ng/ul	299535	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	38
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	38
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	35
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	30
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

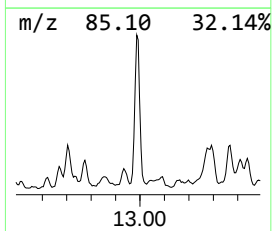
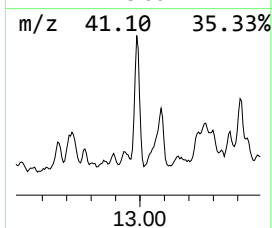
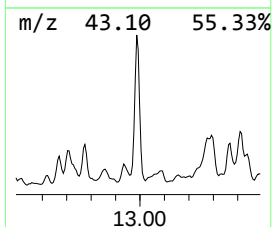
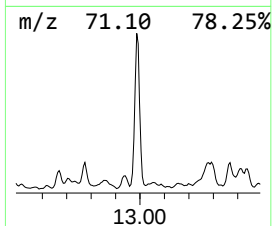
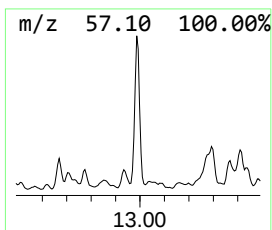
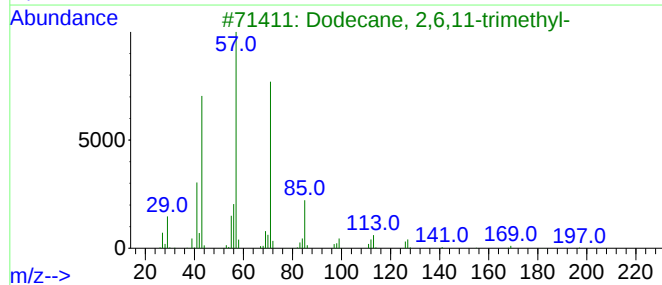
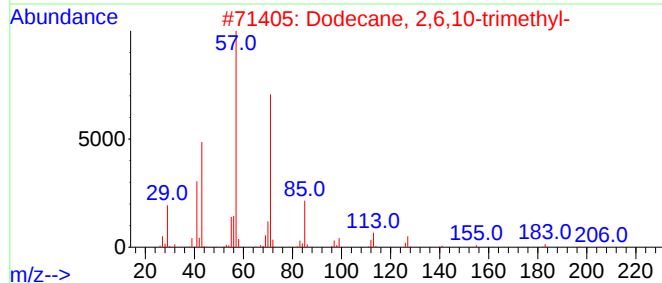
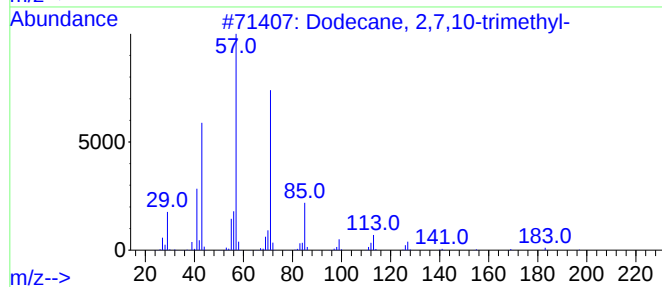
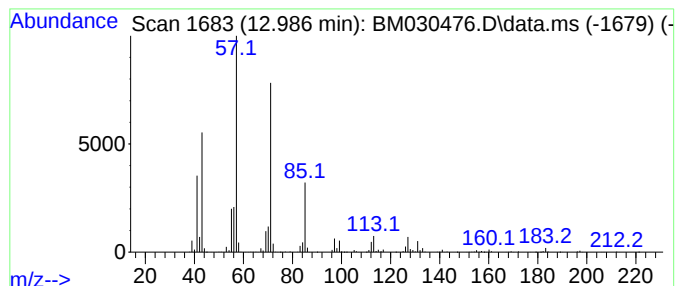
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.990	22.60 ng/ul	2346840	Acenaphthene-d10			14.451
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	91
2	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	90
3	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	83
4	Decane, 5-propyl-		184	C13H28	017312-62-8	74
5	Bacchotricuneatin c		342	C20H22O5	066563-30-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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Misc :
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ClientSampleId :
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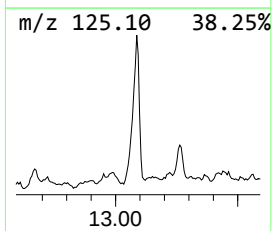
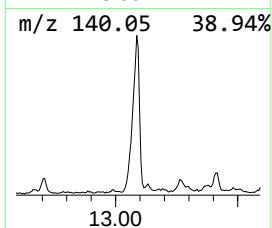
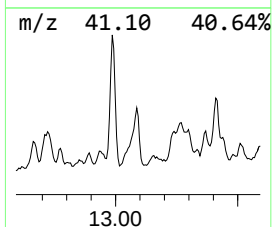
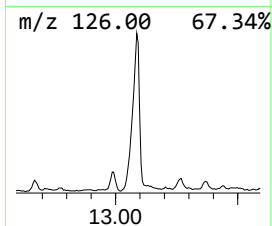
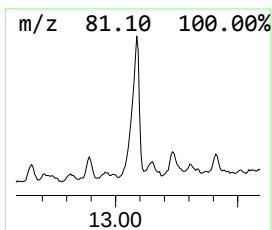
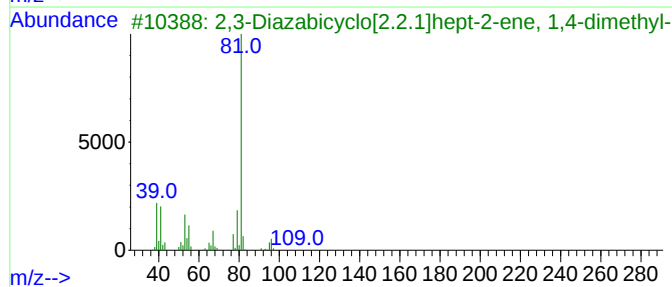
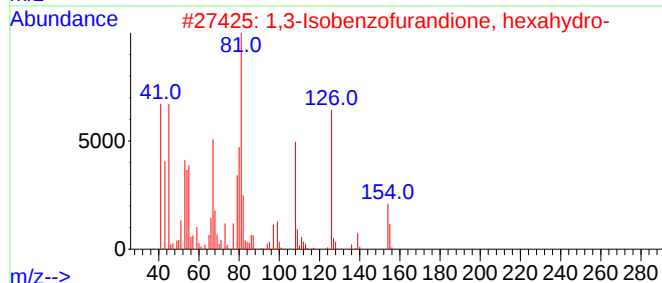
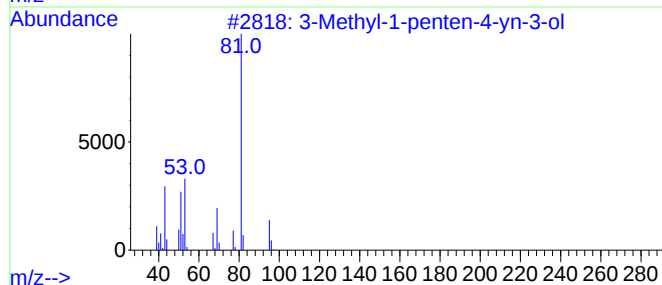
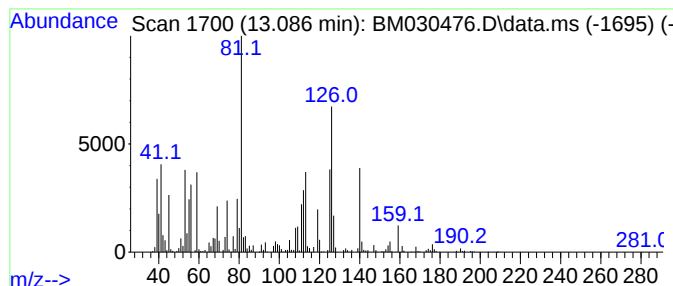
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 unknown-05 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.090	17.17 ng/ul	1782620	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methyl-1-penten-4-yn-3-ol	96	C6H8O	003230-69-1	38
2			1,3-Isobenzofurandione, hexahydro-	154	C8H10O3	000085-42-7	38
3			2,3-Diazabicyclo[2.2.1]hept-2-en...	124	C7H12N2	071312-54-4	32
4			2-Furanpropionic acid	140	C7H8O3	000935-13-7	16
5			3-Methyl-2-furoic acid	126	C6H6O3	004412-96-8	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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Sample : M2596-05
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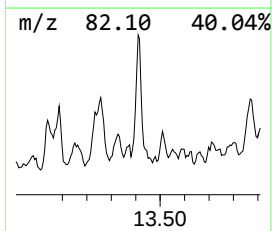
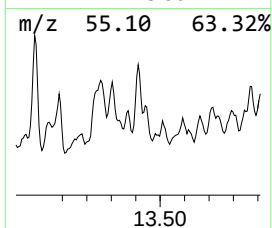
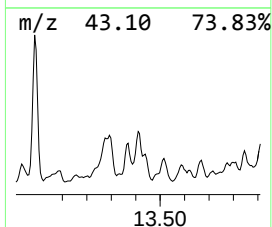
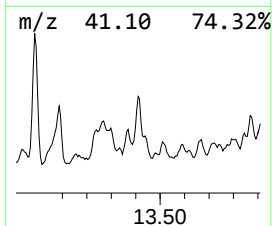
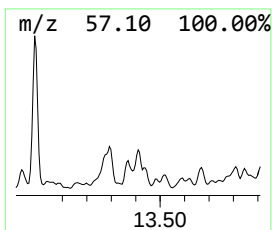
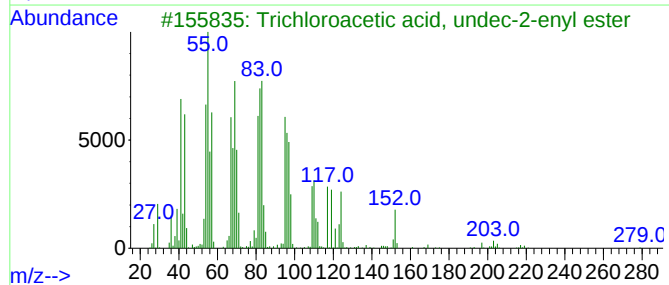
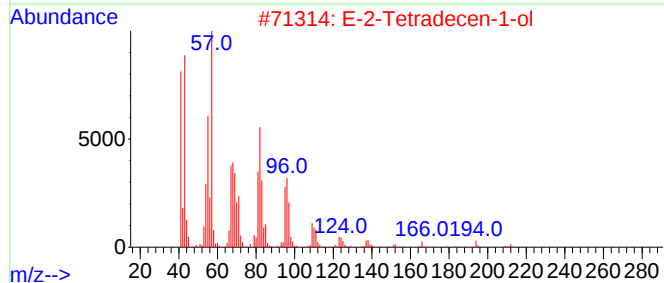
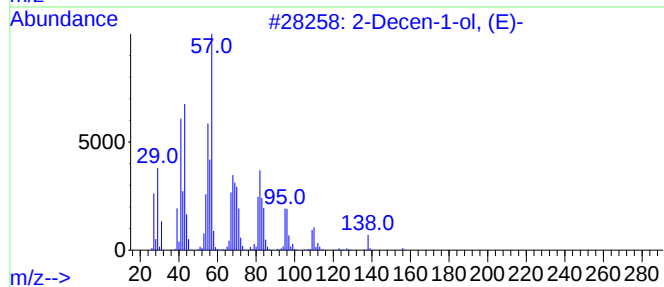
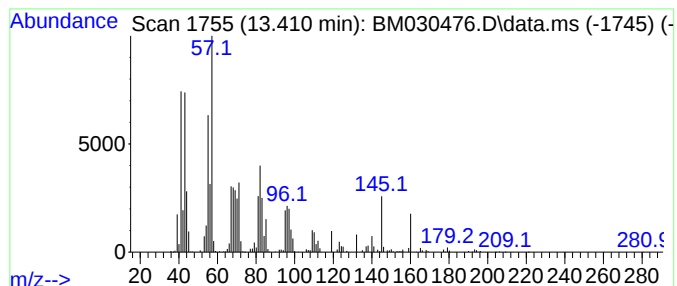
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 2-Decen-1-ol, (E)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.410	15.03 ng/ul	1560610	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Decen-1-ol, (E)-	156	C10H20O	018409-18-2	70
2	E-2-Tetradecen-1-ol	212	C14H28O	1000130-83-7	58
3	Trichloroacetic acid, undec-2-en...	314	C13H21Cl3O2	1000299-26-1	58
4	Dodecanal	184	C12H24O	000112-54-9	58
5	1,14-Tetradecanediol	230	C14H30O2	019812-64-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

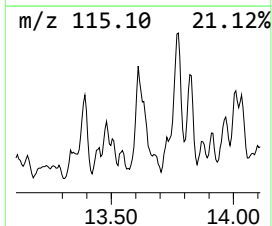
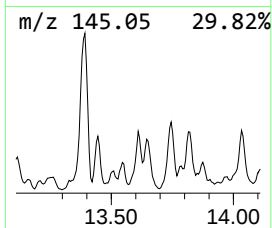
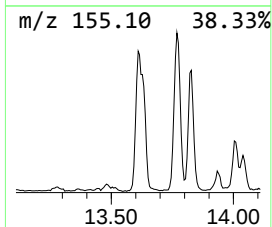
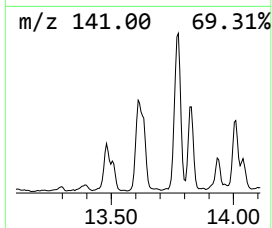
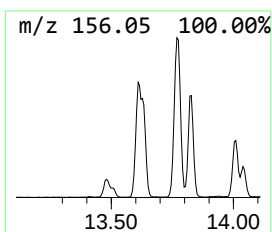
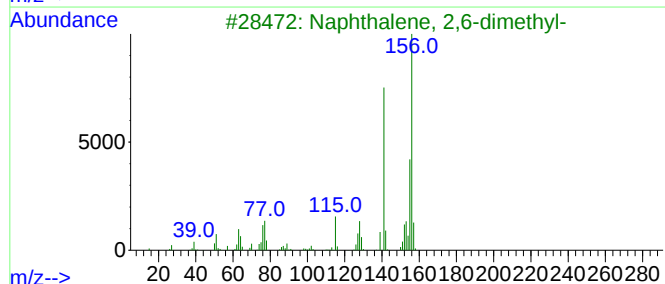
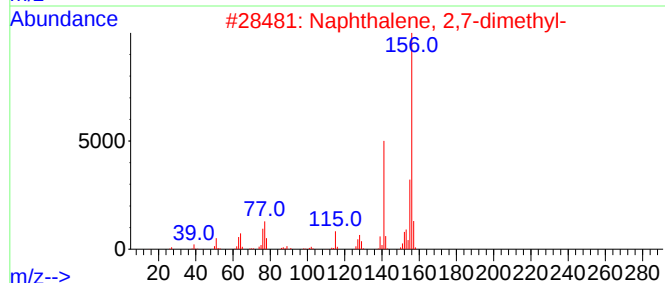
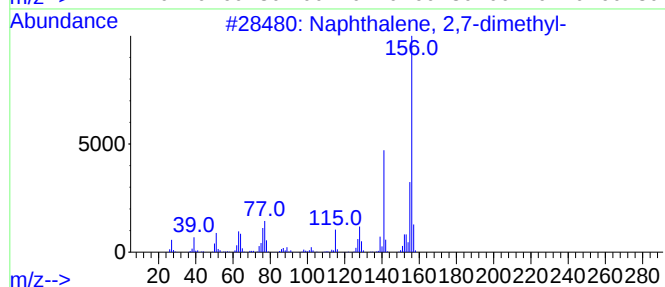
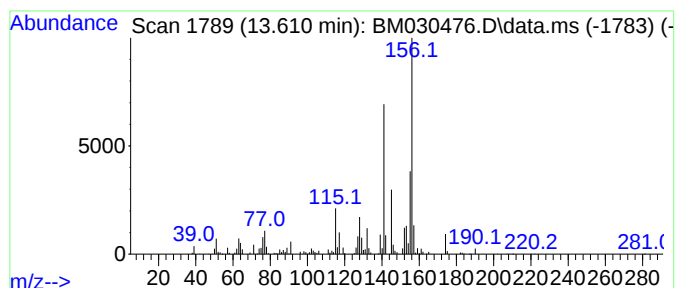
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 Naphthalene, 2,7-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.610	15.51 ng/ul	1610380	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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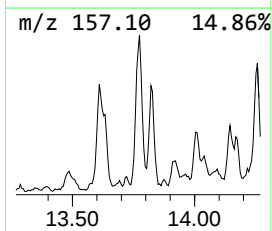
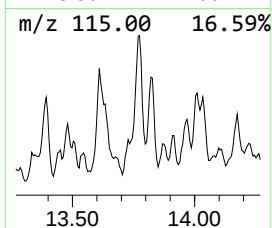
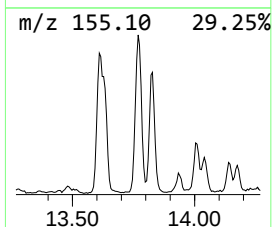
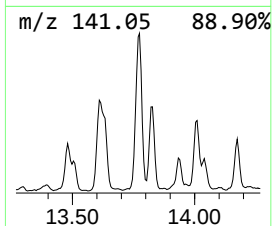
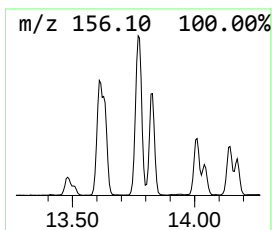
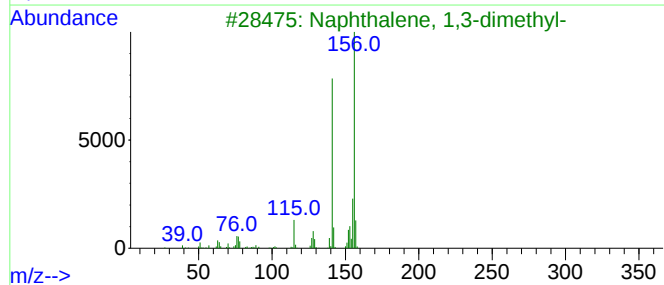
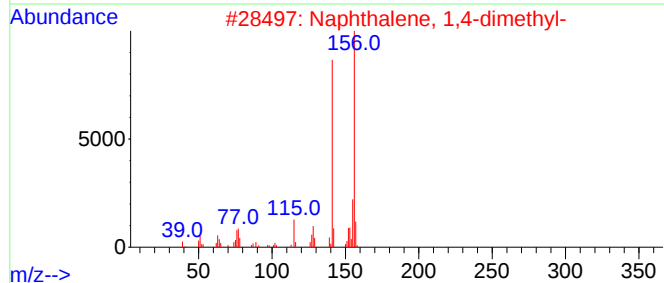
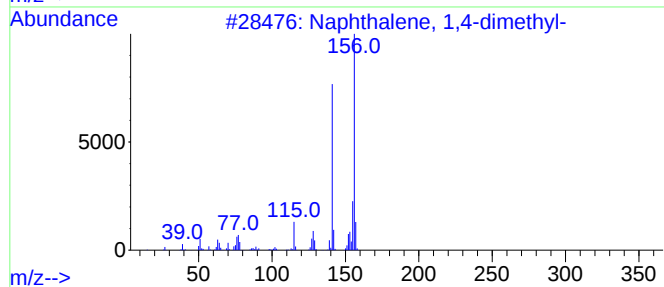
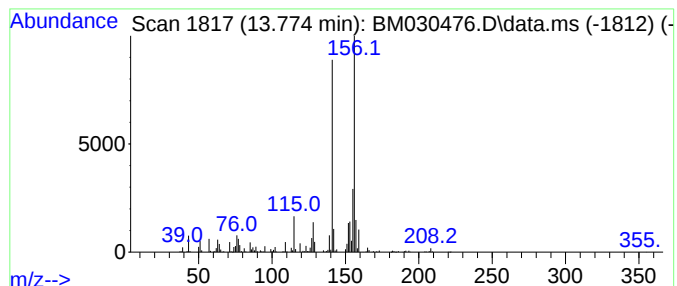
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 Naphthalene, 1,4-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.770	8.95 ng/ul	929513	Acenaphthene-d10	14.451

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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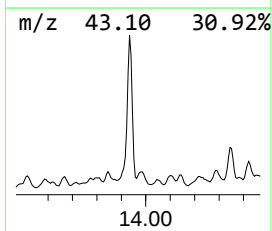
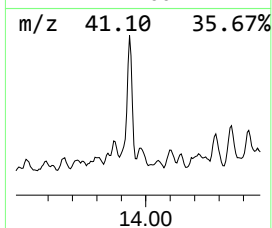
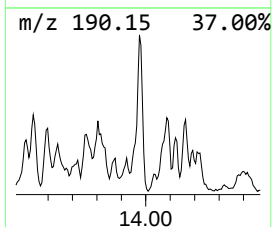
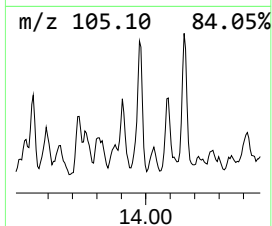
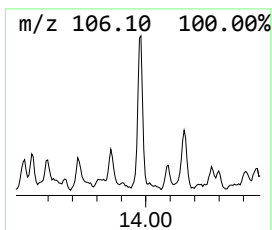
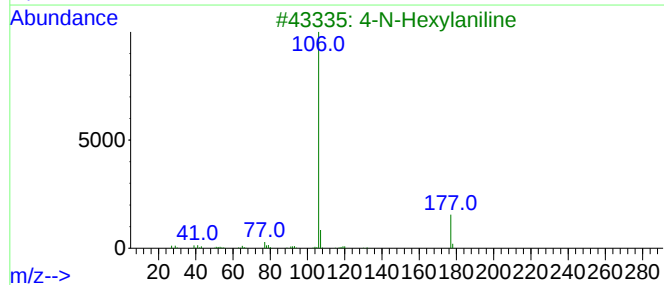
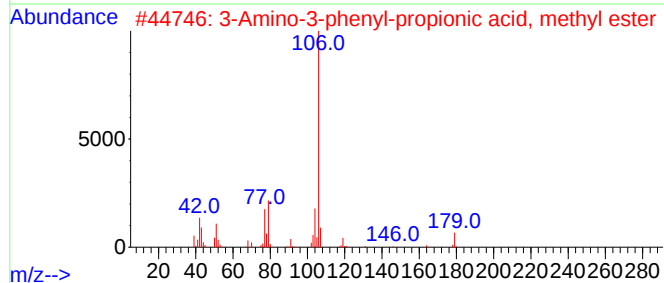
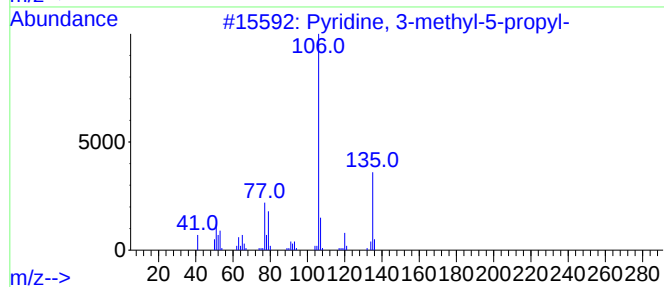
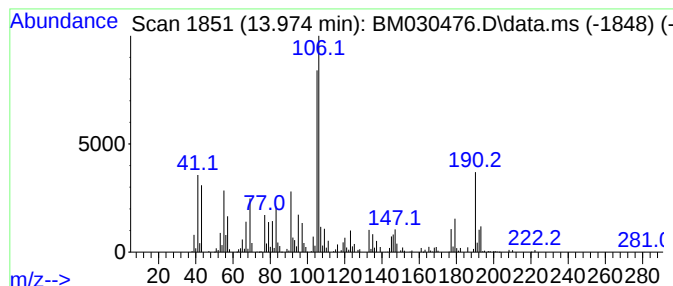
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 unknown-06 Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	3.30 ng/ul	342517	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 3-methyl-5-propyl-	135	C9H13N	053957-26-9	46
2			3-Amino-3-phenyl-propionic acid,...	179	C10H13NO2	1000317-52-0	43
3			4-N-Hexylaniline	177	C12H19N	033228-45-4	43
4			(3-Methylphenyl) methanol, 2-met...	192	C13H20O	1000374-43-9	43
5			4-N-Hexylaniline	177	C12H19N	033228-45-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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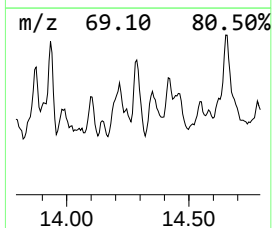
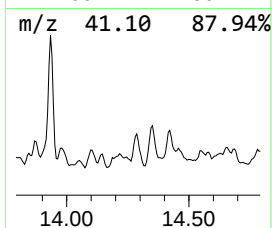
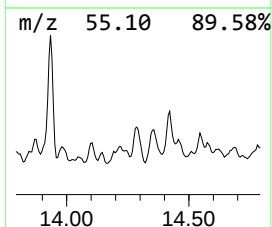
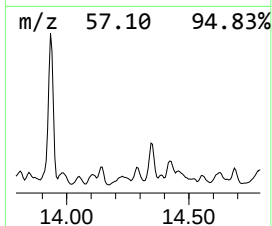
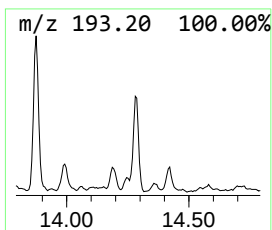
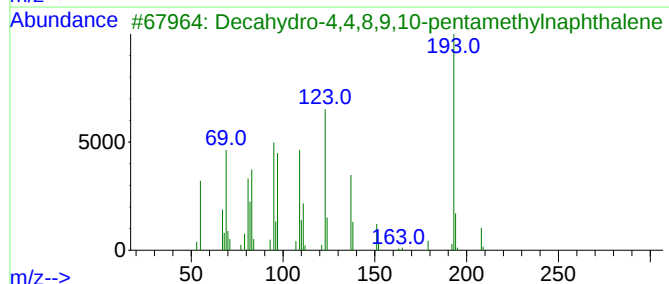
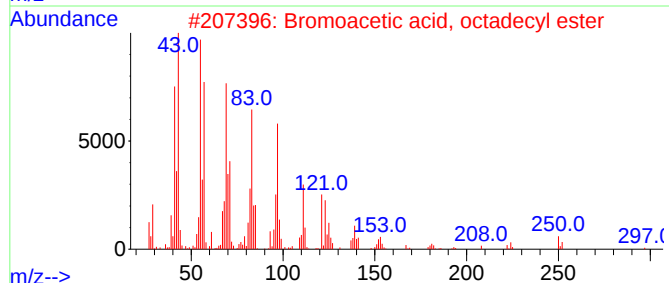
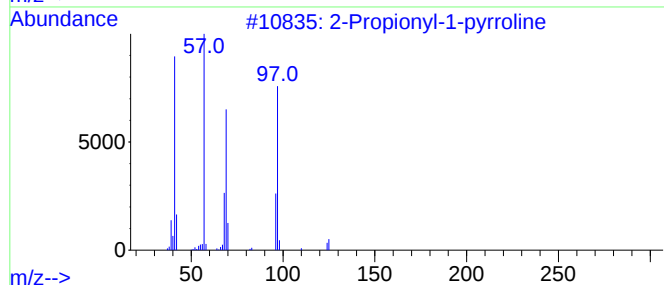
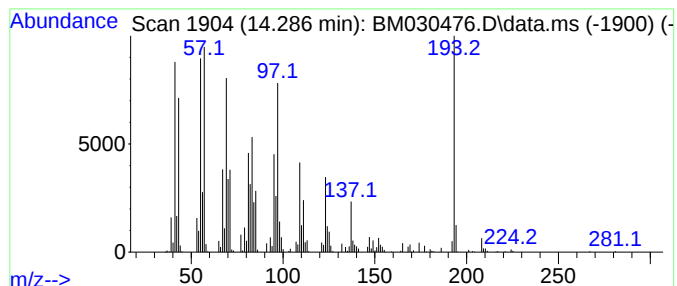
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 unknown-07 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.290	10.95 ng/ul	1136970	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propionyl-1-pyrroline		125	C7H11NO	1000298-55-4	38	
2	Bromoacetic acid, octadecyl ester		390	C20H39BrO2	018992-03-5	38	
3	Decahydro-4,4,8,9,10-pentamethyl...		208	C15H28	080655-44-3	38	
4	1,7-Nonadiene, 4,8-dimethyl-		152	C11H20	062108-28-5	30	
5	Trifluoroacetic acid, pentadecyl...		324	C17H31F3O2	959010-23-2	30	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

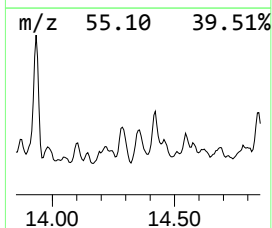
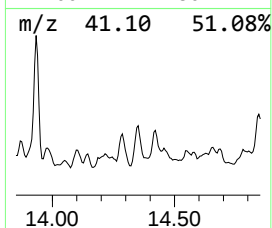
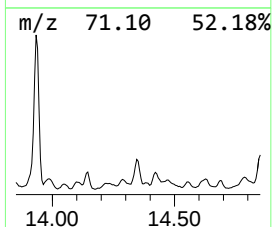
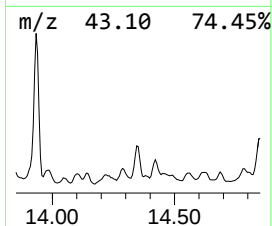
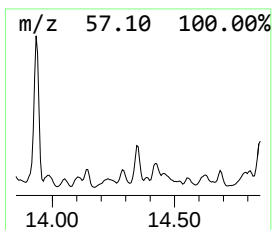
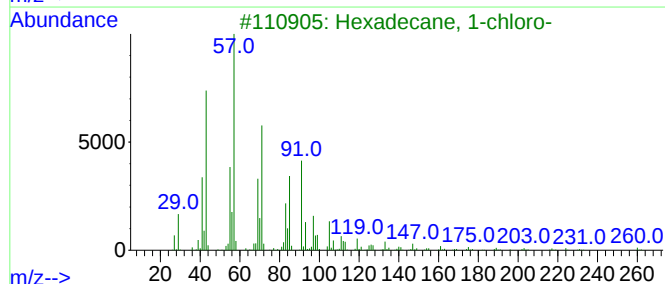
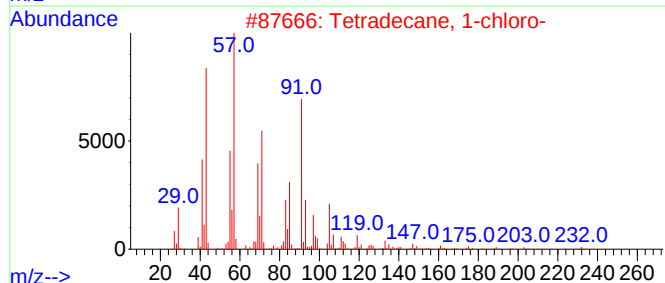
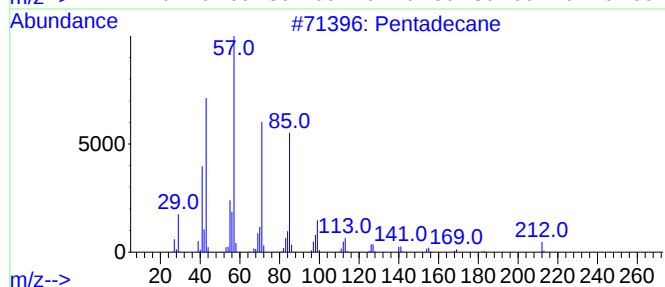
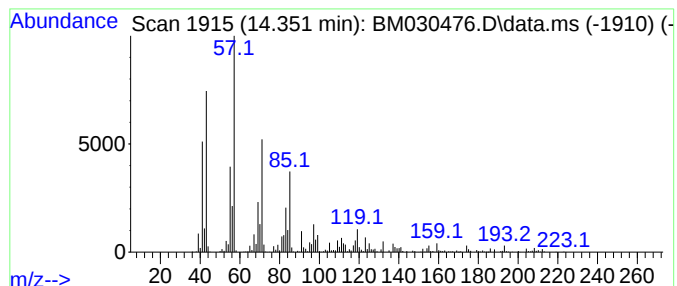
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.350	12.12 ng/ul	1258100	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	93
2	Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	72
3	Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	72
4	Dodecane	170	C12H26	000112-40-3	64
5	Pentadecane	212	C15H32	000629-62-9	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

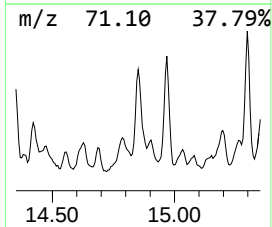
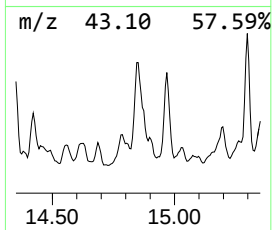
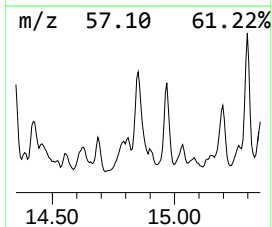
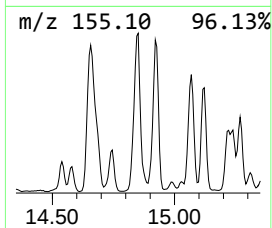
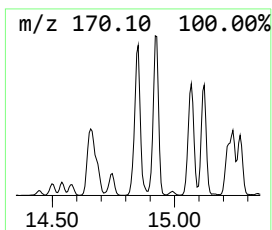
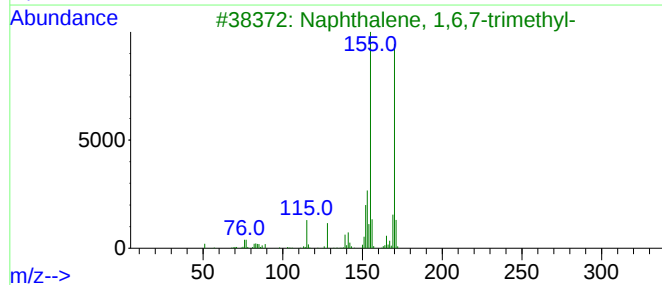
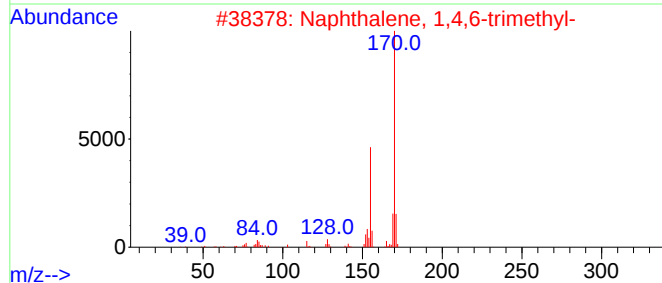
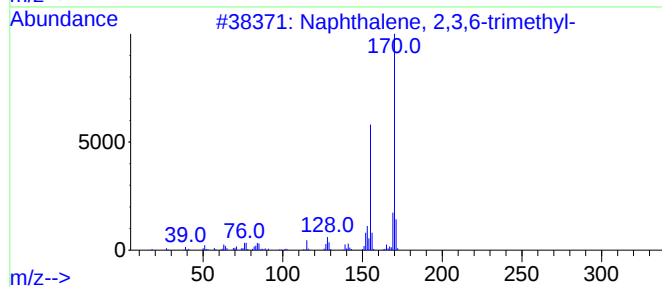
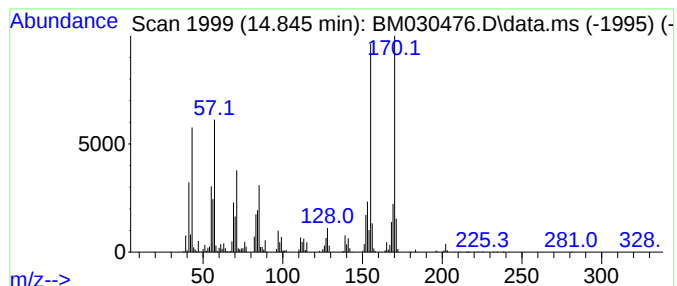
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 Naphthalene, 2,3,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.850	24.36 ng/ul	2529410	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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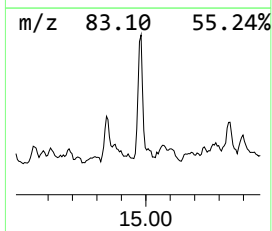
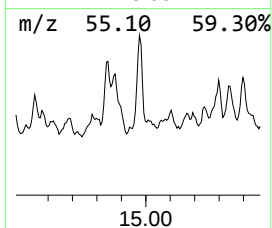
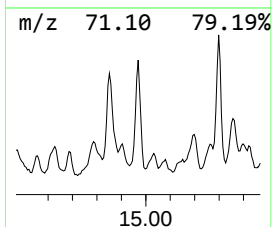
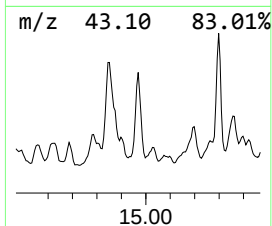
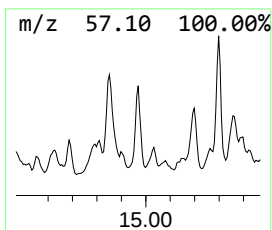
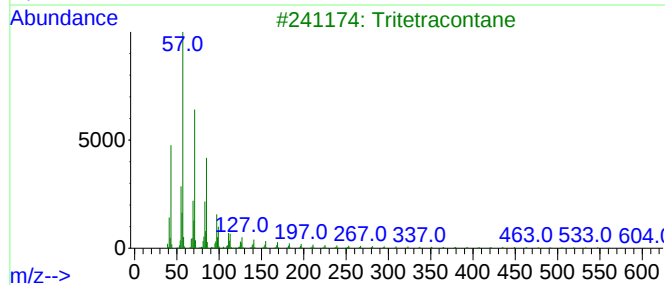
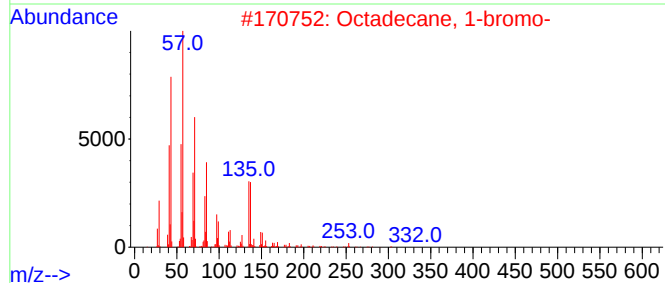
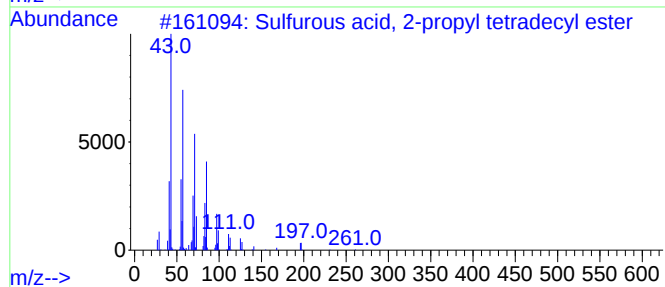
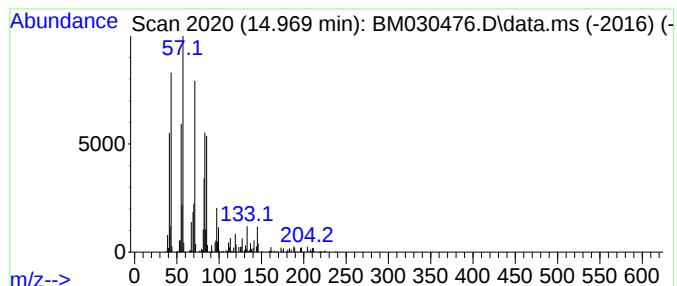
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 Sulfurous acid, 2-propyl te... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.970	16.39 ng/ul	1701920	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-propyl tetradecyl ester	320	C17H36O3S	1000309-12-5	58
2			Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	58
3			Tritetracontane	605	C43H88	007098-21-7	58
4			Nonadecane	268	C19H40	000629-92-5	53
5			Tetratetracontane	619	C44H90	007098-22-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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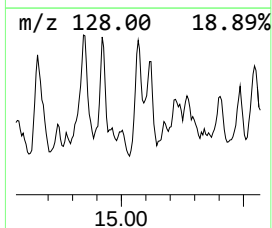
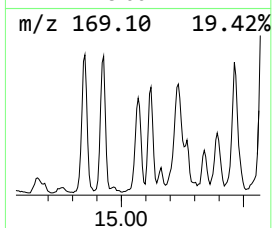
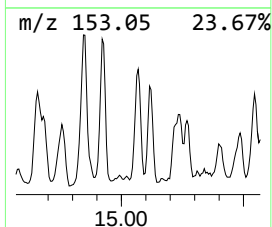
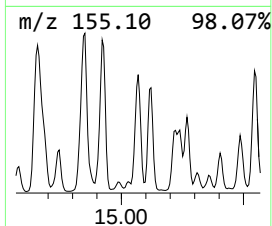
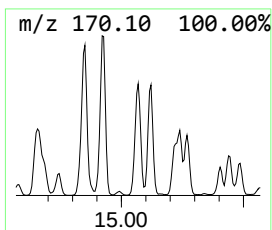
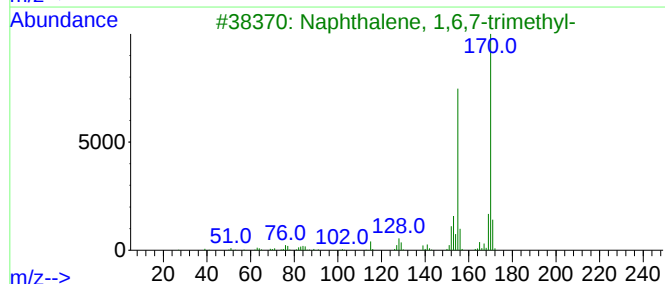
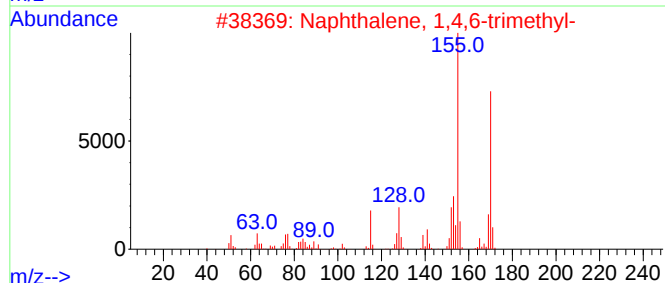
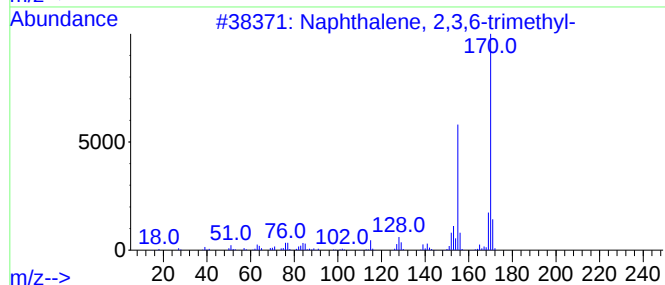
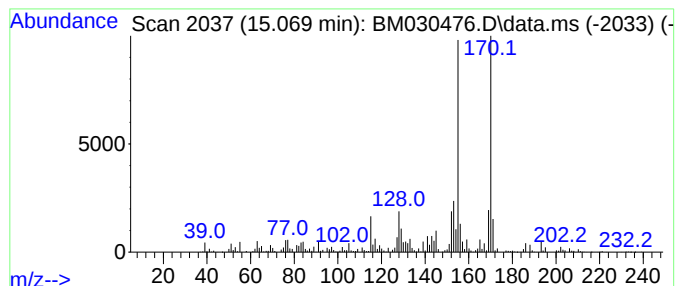
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 Naphthalene, 1,4,6-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.070	11.70 ng/ul	1214590	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
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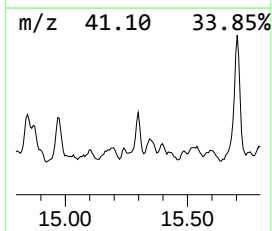
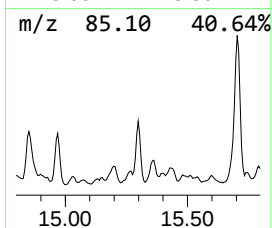
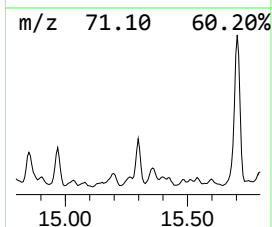
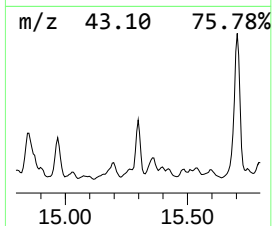
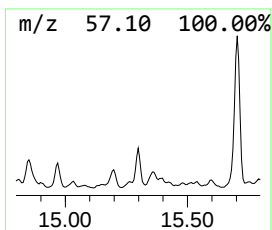
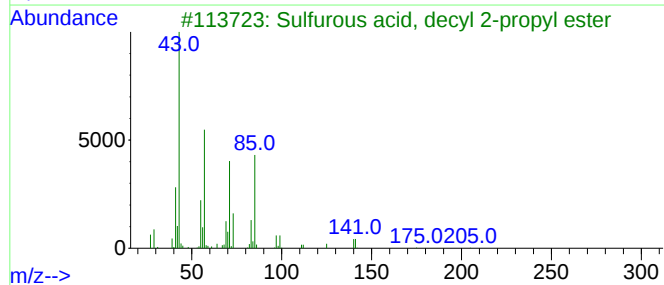
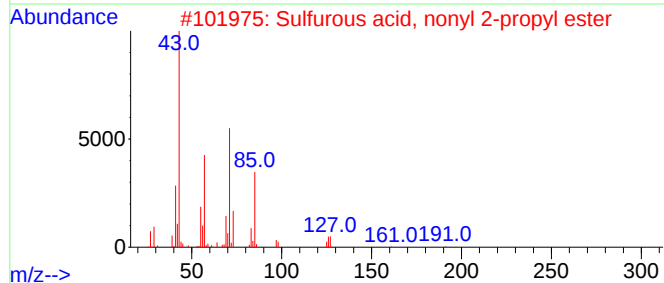
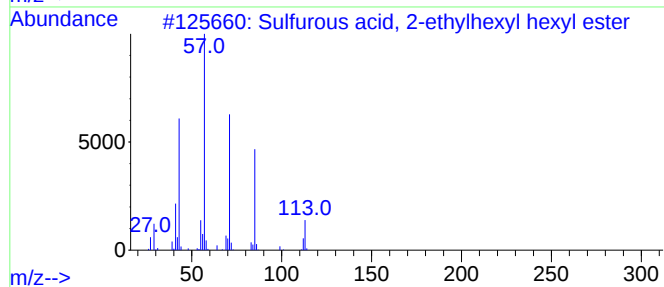
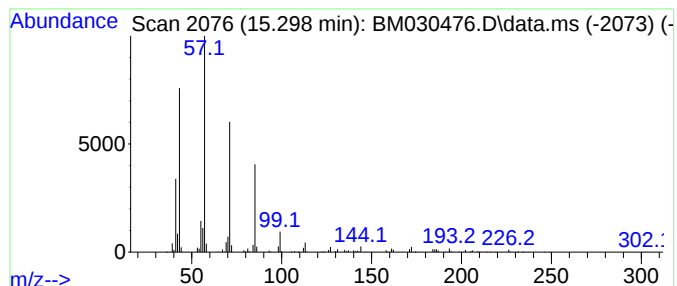
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 Sulfurous acid, 2-ethylhexy... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.300	8.12 ng/ul	843329	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	72
2			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	64
3			Sulfurous acid, decyl 2-propyl e...	264	C13H28O3S	1000309-12-1	64
4			Hexadecane	226	C16H34	000544-76-3	49
5			Octadecane, 5-methyl-	268	C19H40	025117-35-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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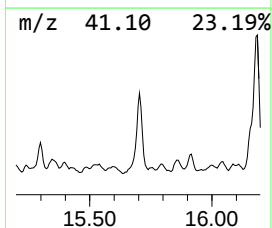
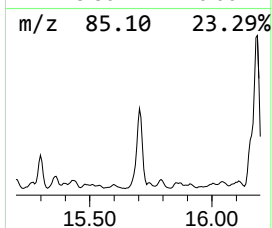
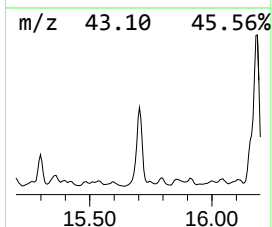
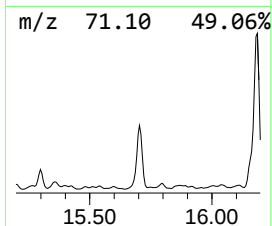
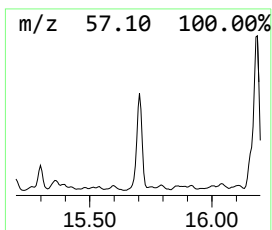
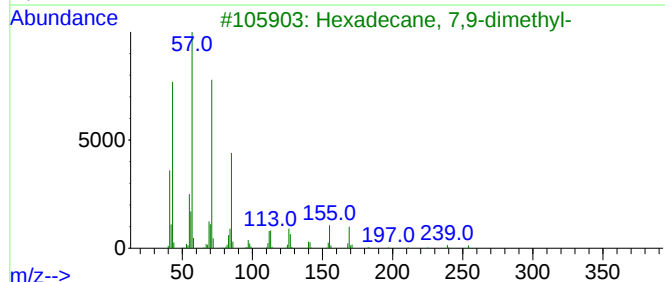
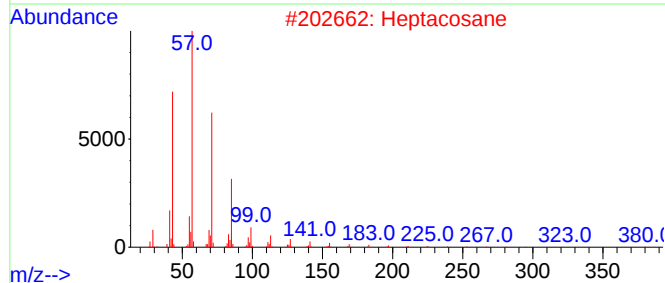
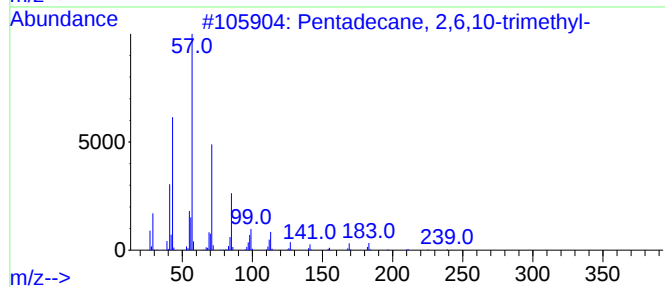
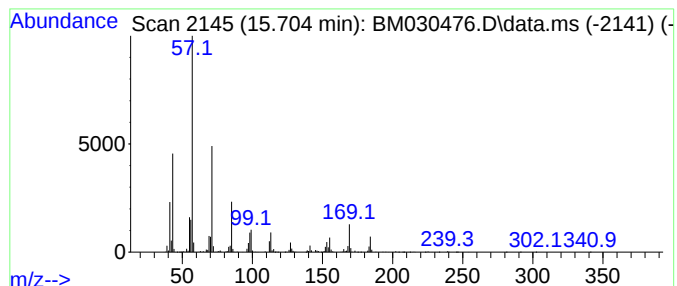
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.700	40.12 ng/ul	4166100	Acenaphthene-d10	14.451

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Heptacosane	380	C27H56	000593-49-7	72
3		Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	58
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	53
5		Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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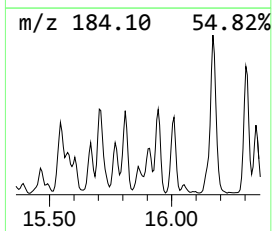
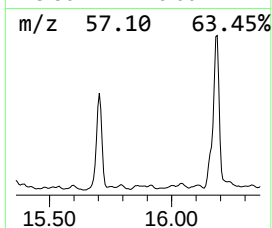
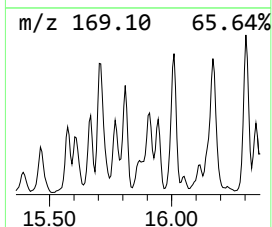
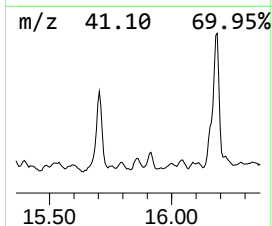
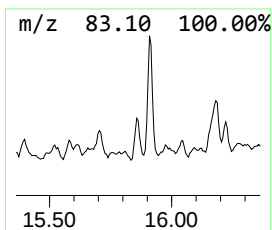
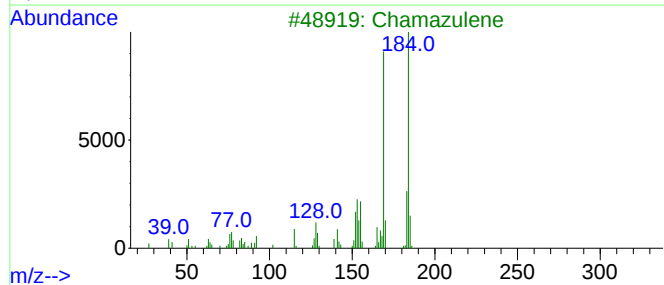
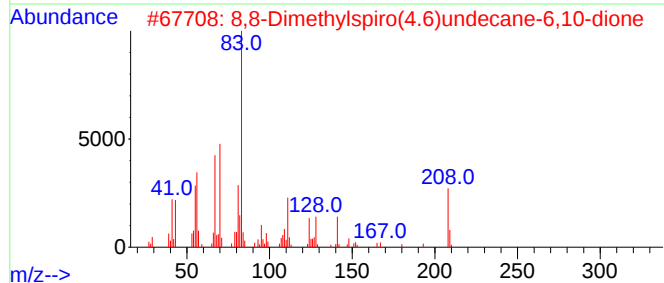
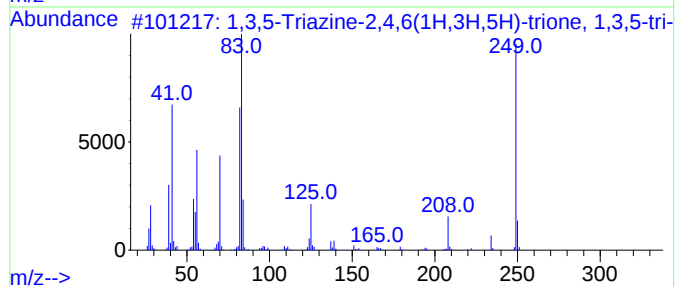
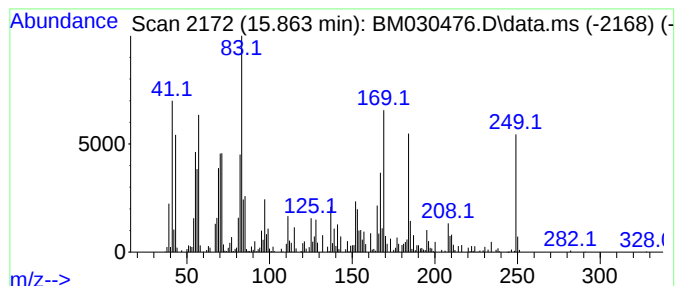
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 unknown-08 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.860	5.68 ng/ul	706998	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	49
2			8,8-Dimethylspiro(4.6)undecane-6...	208	C13H20O2	068181-81-7	35
3			Chamazulene	184	C14H16	000529-05-5	30
4			[1,1'-Biphenyl]-4-methanol	184	C13H12O	003597-91-9	25
5			1-Acenaphthylenol, 1,2-dihydro-1...	184	C13H12O	041002-74-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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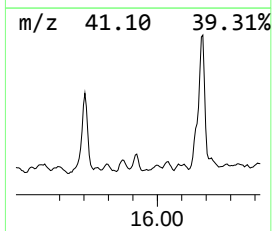
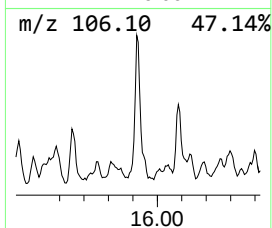
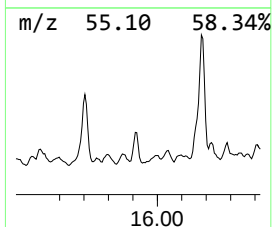
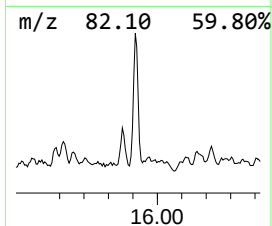
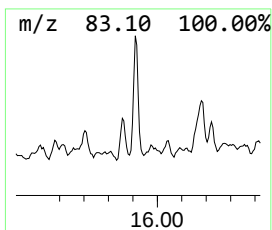
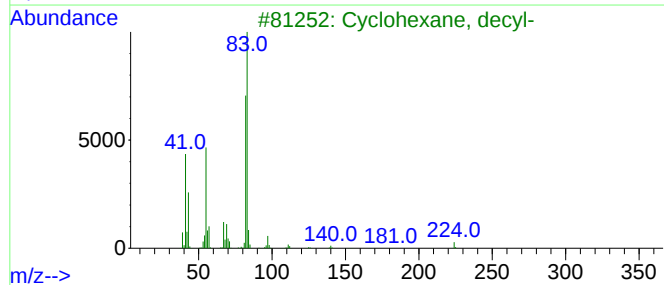
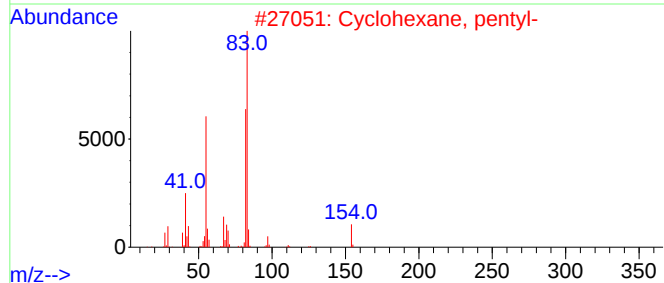
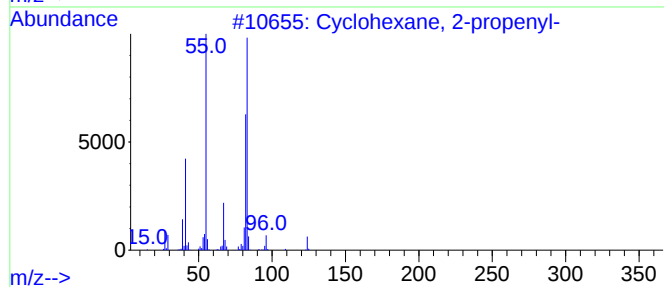
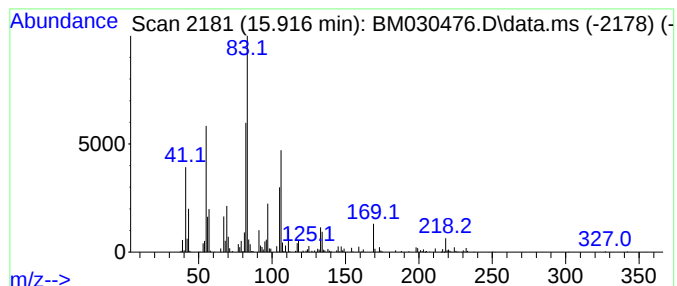
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 unknown-09 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.920	5.23 ng/ul	650910	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	45
2			Cyclohexane, pentyl-	154	C11H22	004292-92-6	43
3			Cyclohexane, decyl-	224	C16H32	001795-16-0	43
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	38
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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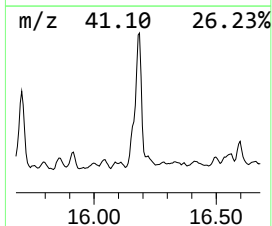
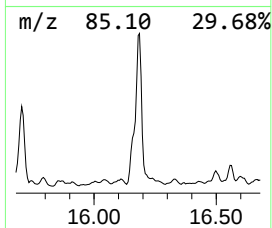
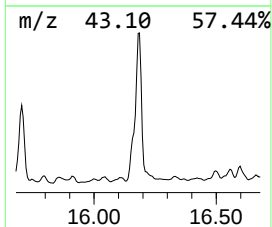
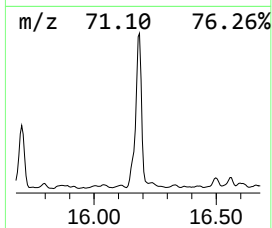
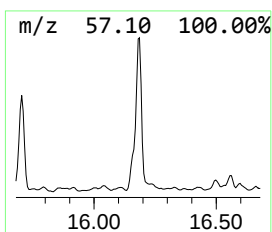
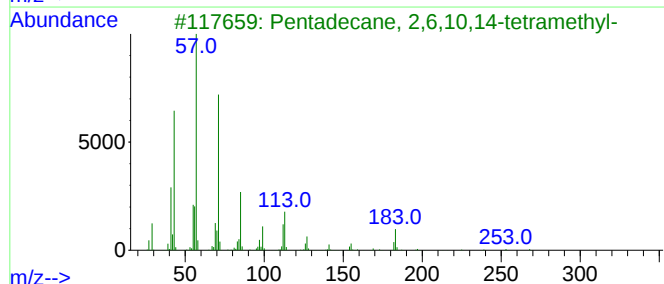
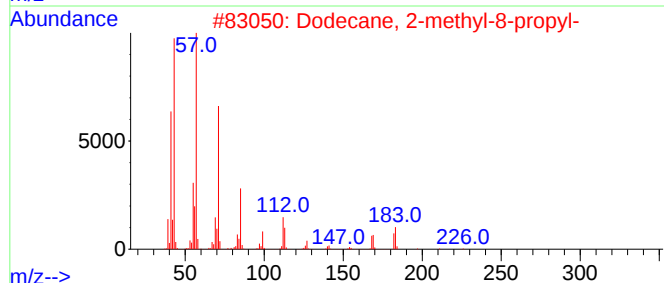
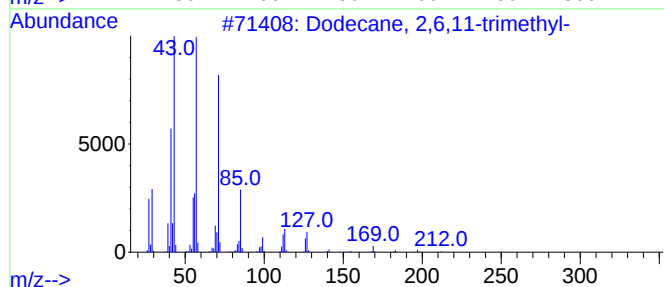
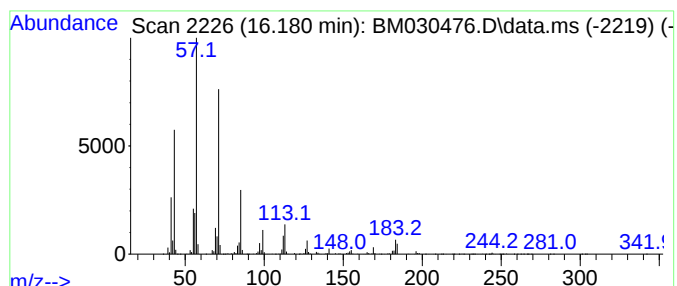
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.180	77.48 ng/ul	9646650	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	94
2	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	94
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
4	Tridecane, 5-propyl-	226	C16H34	055045-11-9	87
5	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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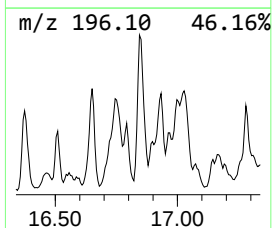
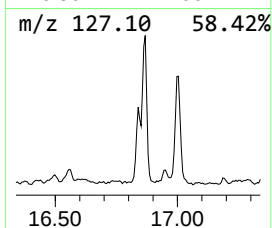
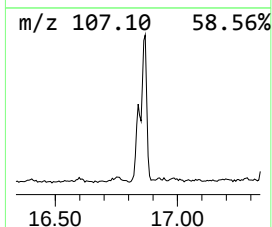
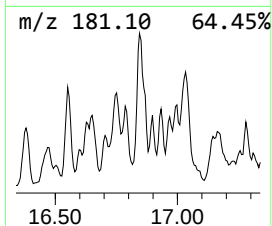
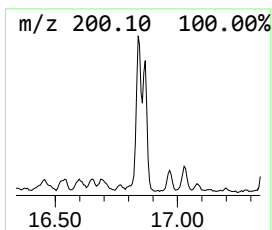
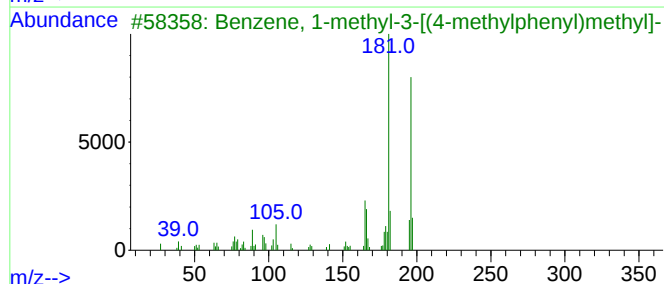
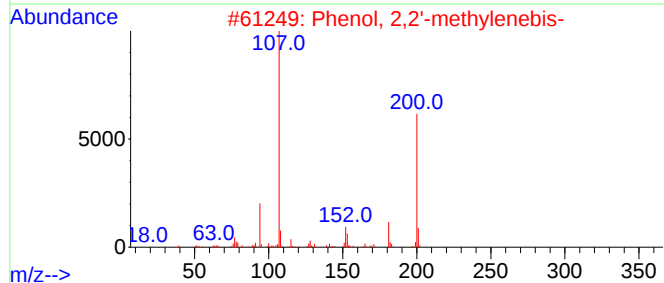
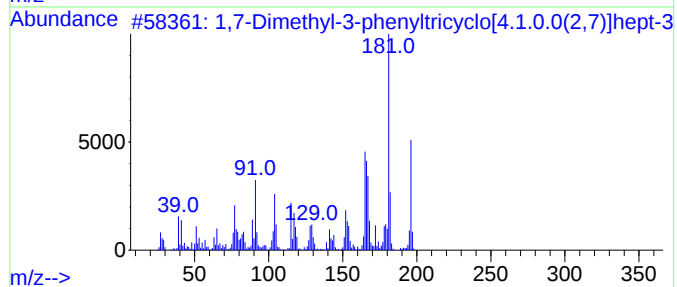
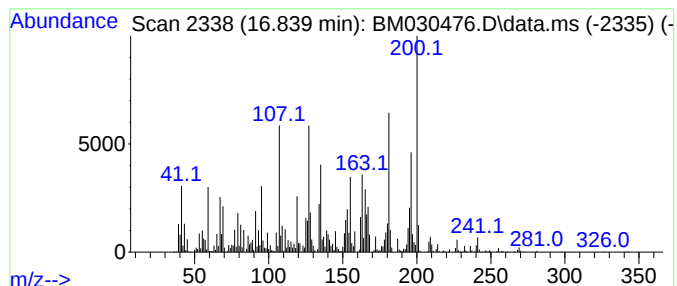
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 unknown-10 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.840	10.56 ng/ul	1314330	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,7-Dimethyl-3-phenyltricyclo[4.1.0.0(2,7)]hept-3	196	C15H16	1000200-99-4	38
2			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	38
3			Benzene, 1-methyl-3-[(4-methylphenyl)methyl]-	196	C15H16	021895-16-9	25
4			Methane, di-p-tolyl-	196	C15H16	004957-14-6	25
5			Benzene, 2,6-dimethyl-1-(phenylmethyl)-	196	C15H16	028122-29-4	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5Q1

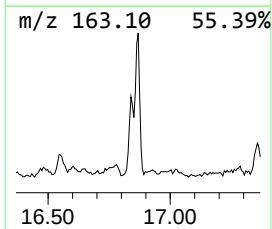
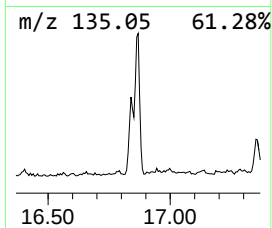
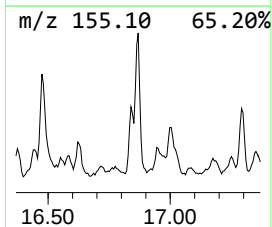
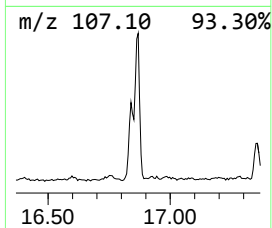
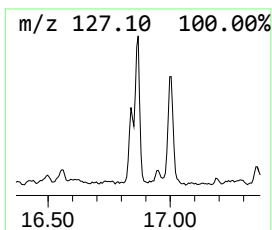
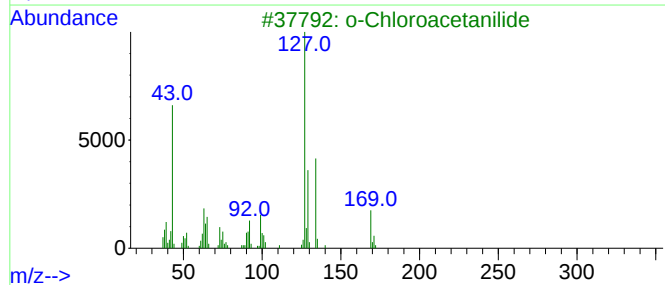
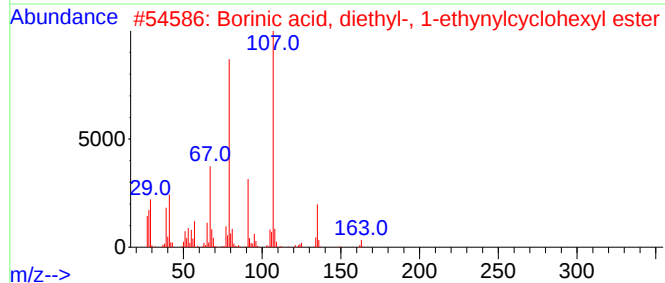
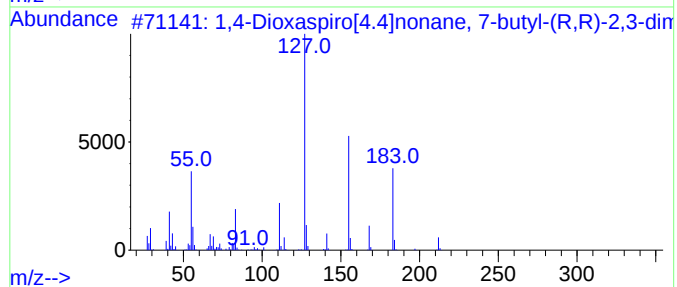
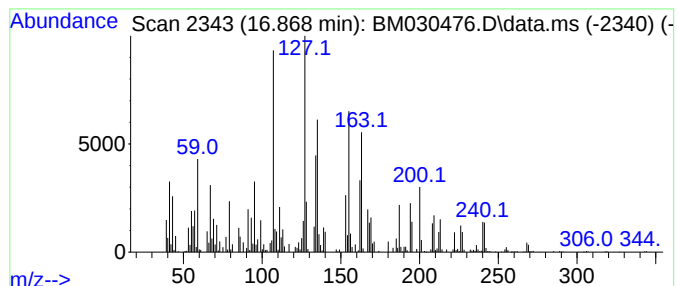
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 unknown-11 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.870	14.60 ng/ul	1818120	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dioxaspiro[4.4]nonane, 7-but...	212	C13H24O2	104807-77-4	18
2			Borinic acid, diethyl-, 1-ethyny...	192	C12H21BO	055848-34-5	14
3			o-Chloroacetanilide	169	C8H8ClNO	000533-17-5	14
4			Dimethyl (3S,4R,5S,6R)-3,6-dimet...	226	C12H18O4	036337-44-7	14
5			1-Naphthamide, N-allyl-	211	C14H13NO	1000340-22-4	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

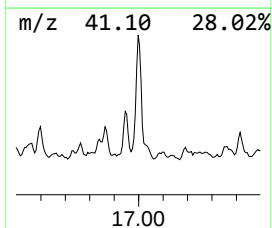
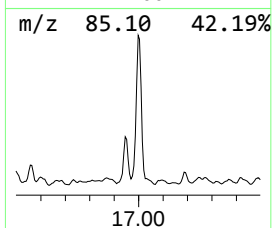
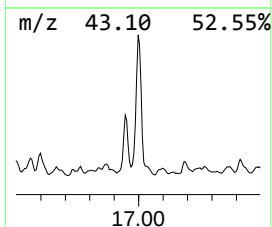
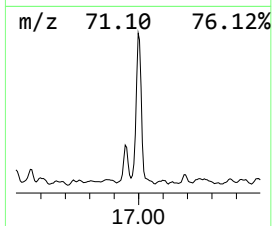
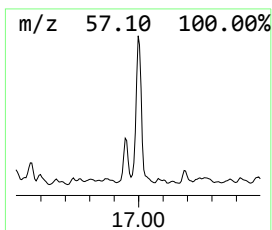
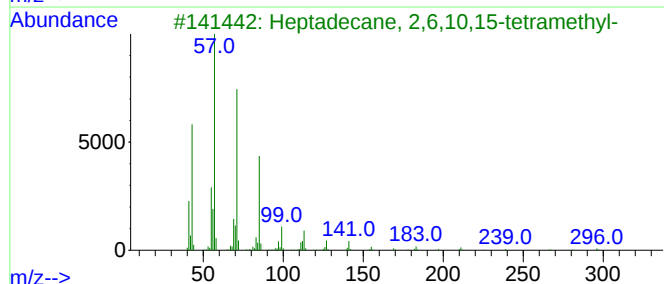
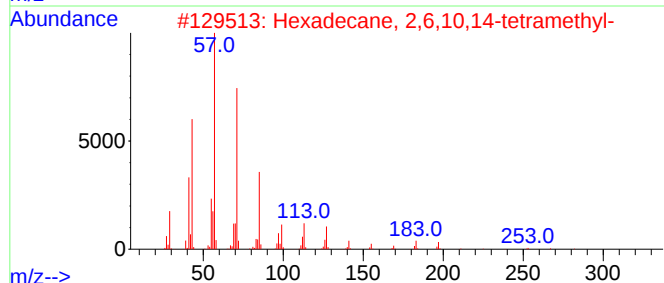
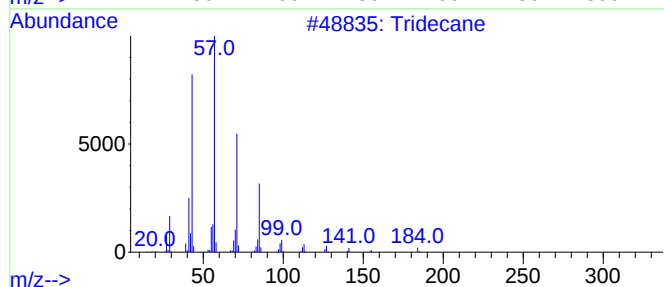
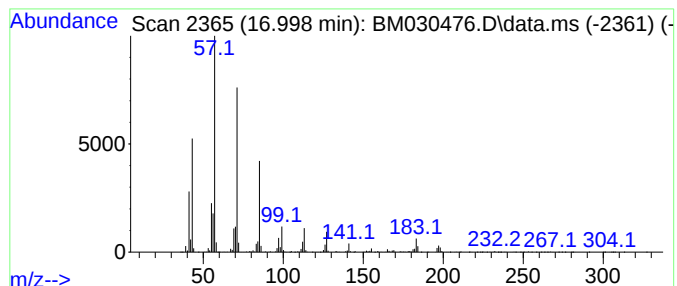
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.000	35.51 ng/ul	4421380	Phenanthrene-d10		17.186	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	91
2	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	91
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	87
4	Eicosane		282	C20H42	000112-95-8	86
5	Pentadecane, 4-methyl-		226	C16H34	002801-87-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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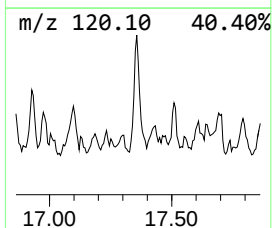
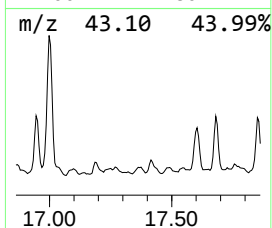
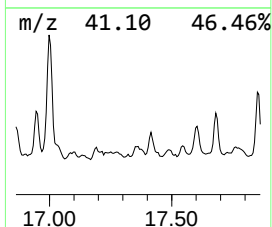
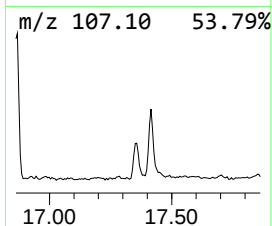
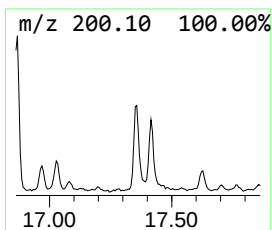
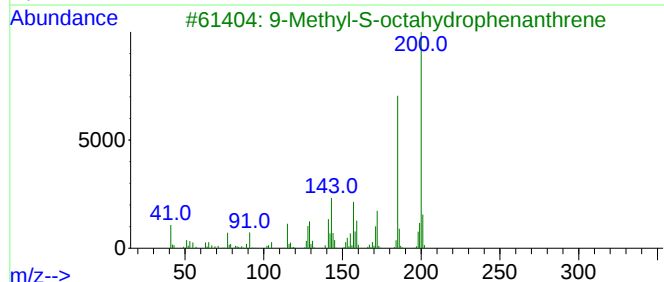
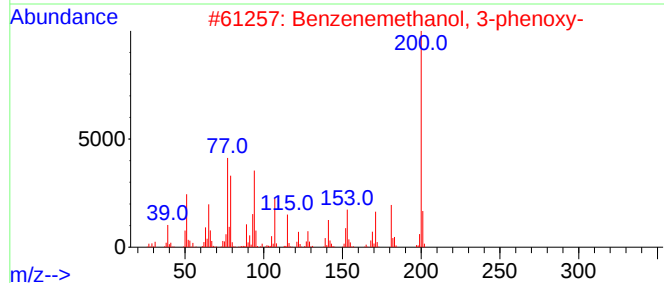
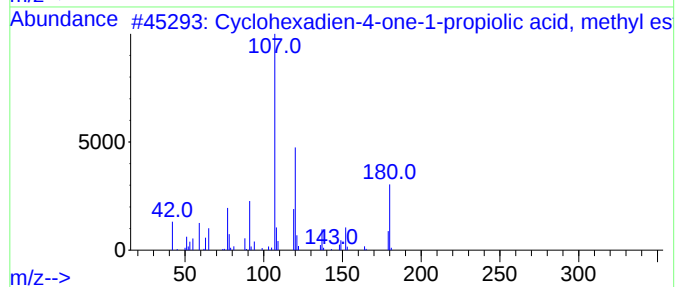
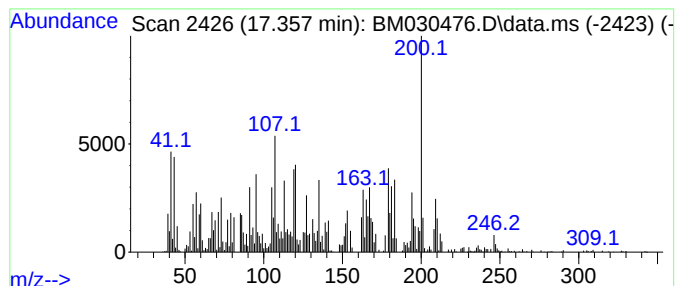
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 unknown-12 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.360	5.67 ng/ul	705924	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadien-4-one-1-propionic ...	180	C10H12O3	1000124-47-3	25
2			Benzenemethanol, 3-phenoxy-	200	C13H12O2	013826-35-2	20
3			9-Methyl-S-octahydrophenanthrene	200	C15H20	023861-81-6	20
4			Benzoic acid, 2-fluoro-5,6-dimet...	200	C9H9FO4	265670-72-2	18
5			1-(4-Ethoxyphenyl)-1-hydroxy-2-p...	194	C11H14O3	1000123-93-4	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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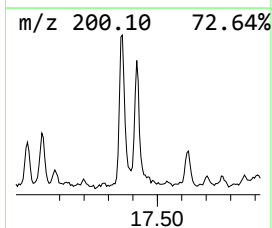
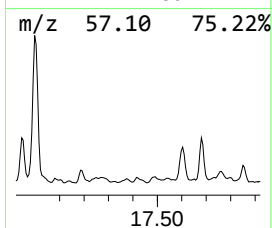
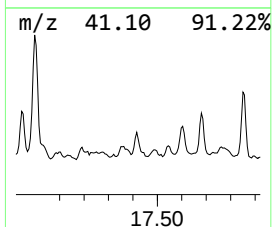
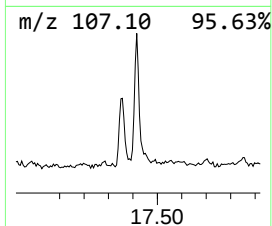
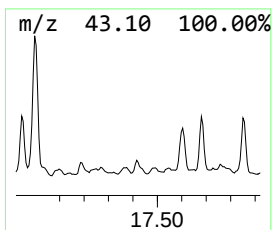
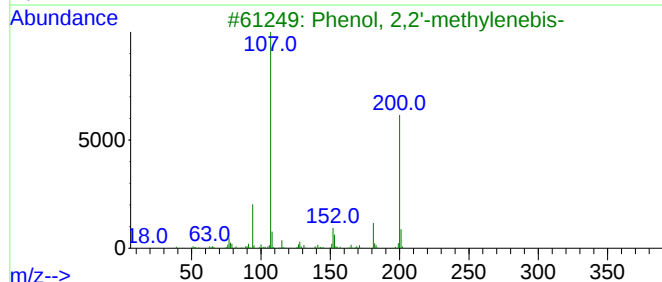
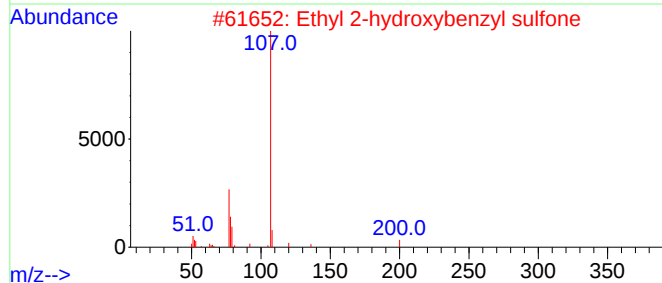
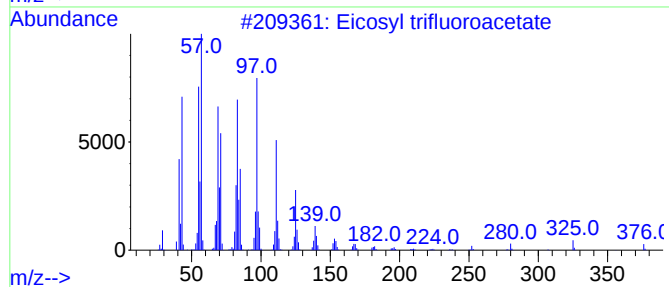
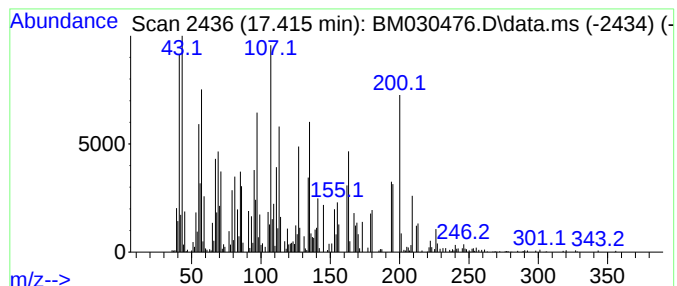
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 unknown-13 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.420	8.58 ng/ul	1068790	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	15
2			Ethyl 2-hydroxybenzyl sulfone	200	C9H12O3S	053380-27-1	14
3			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	14
4			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	14
5			2,4-Pentanedione, 3-(1-methyl-2-...	154	C9H14O2	029149-83-5	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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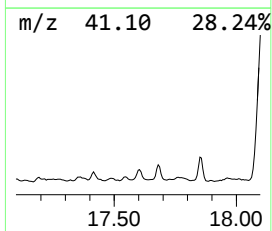
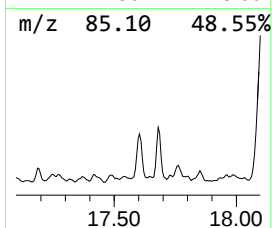
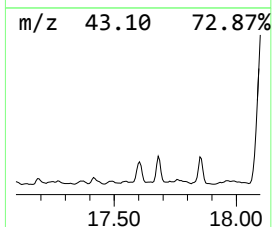
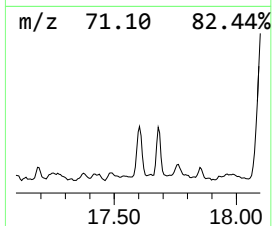
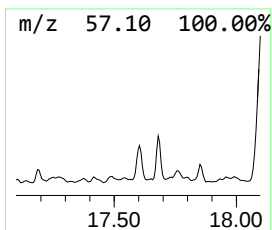
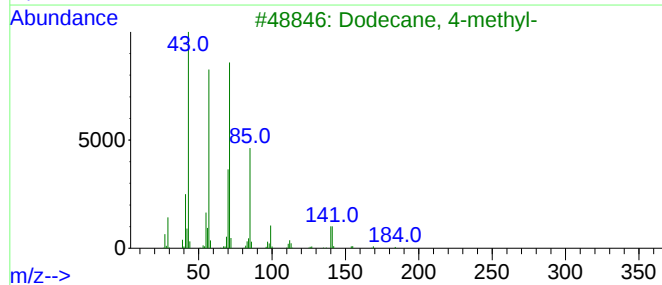
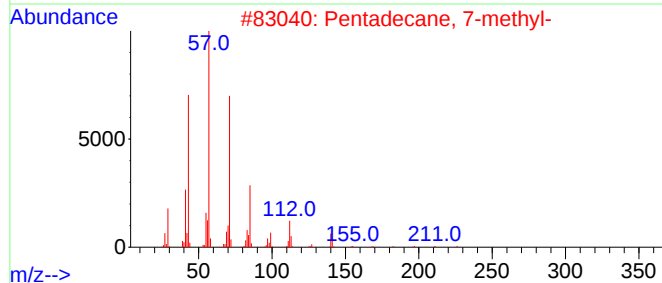
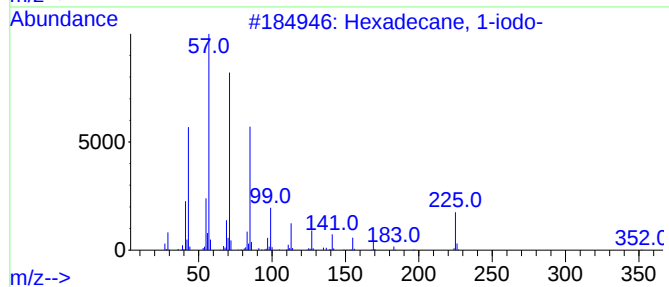
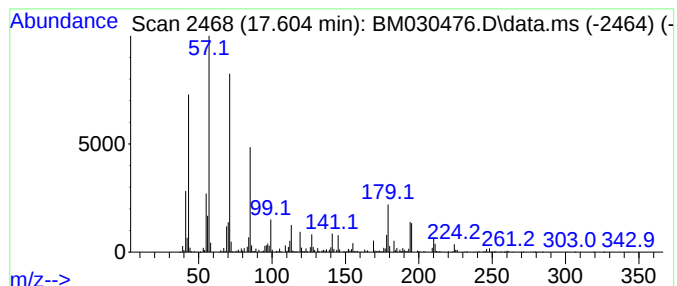
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 Hexadecane, 1-iodo- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.600	15.39 ng/ul	1916070	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	70
2			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	70
3			Dodecane, 4-methyl-	184	C13H28	006117-97-1	64
4			Heptadecane	240	C17H36	000629-78-7	62
5			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	60



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Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
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Instrument :
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ClientSampleId :
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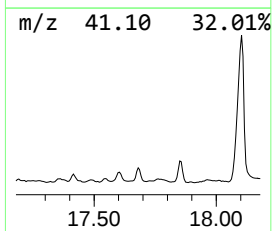
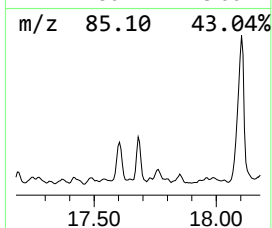
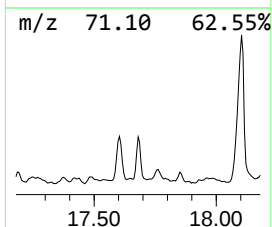
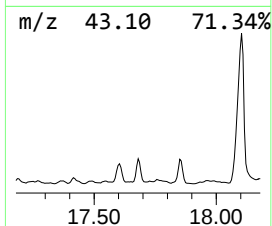
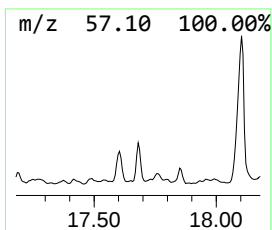
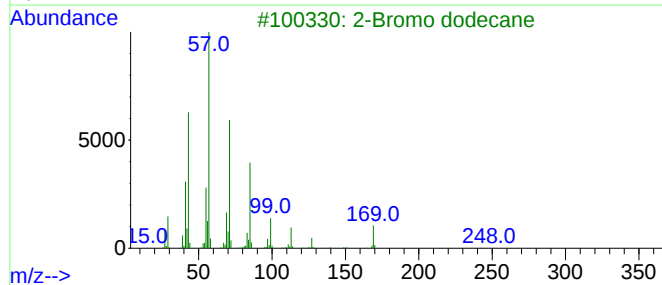
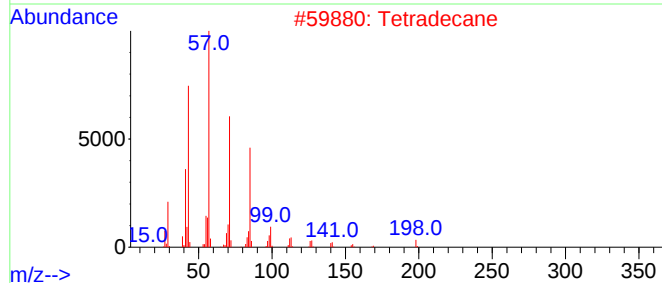
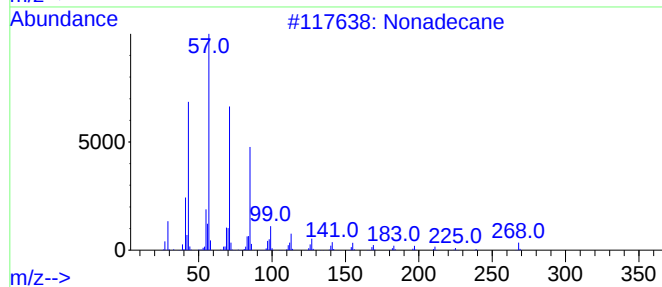
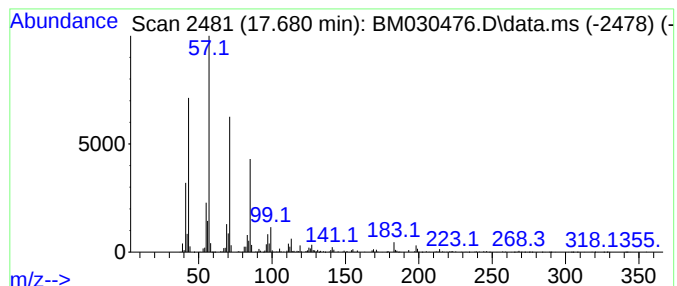
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.680	11.05 ng/ul	1375490	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	95
2			Tetradecane	198	C14H30	000629-59-4	93
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	90
4			Heptadecane	240	C17H36	000629-78-7	86
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
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Instrument :
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ClientSampleId :
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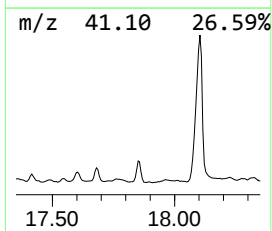
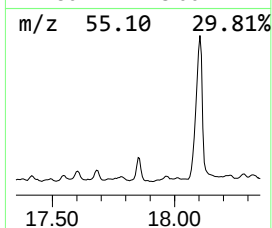
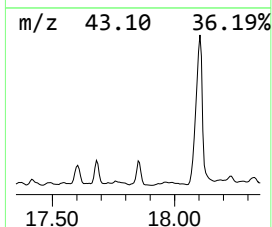
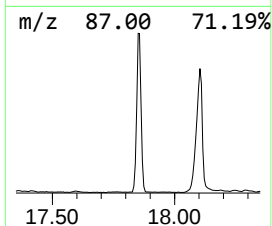
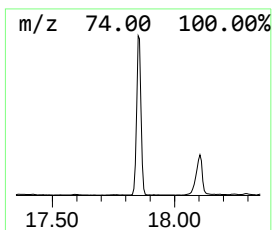
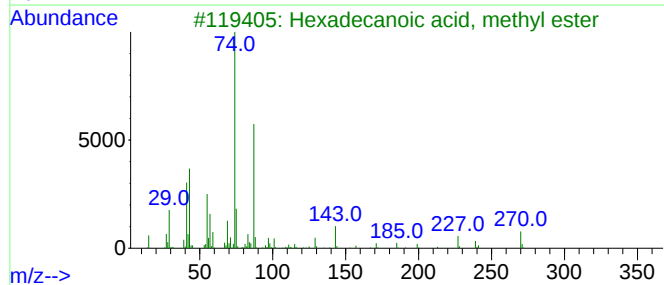
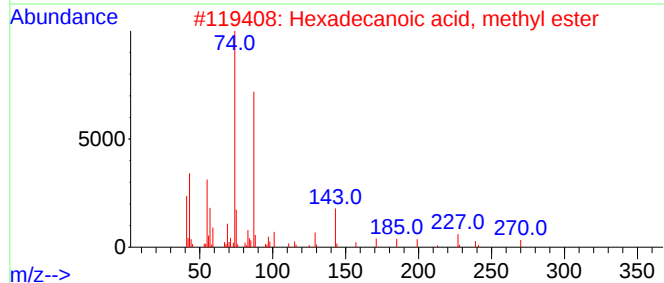
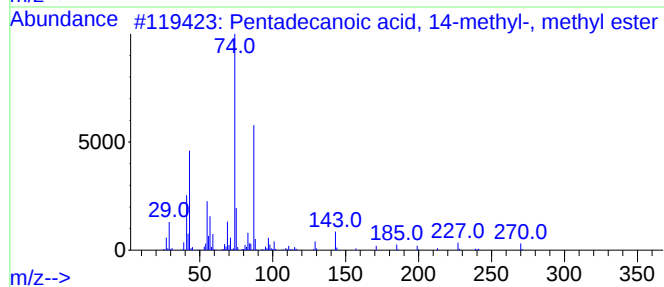
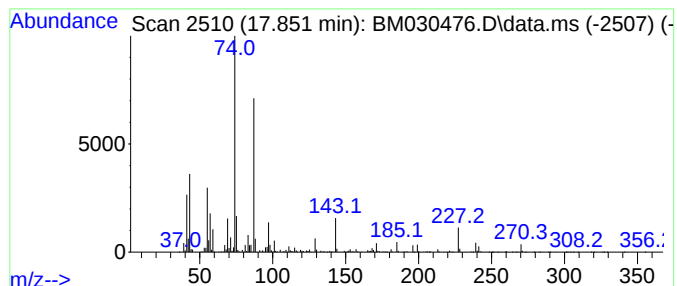
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 Pentadecanoic acid, 14-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	21.63 ng/ul	2693430	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	86
4			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	83
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
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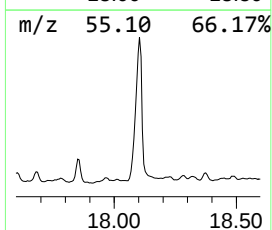
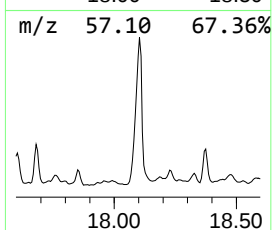
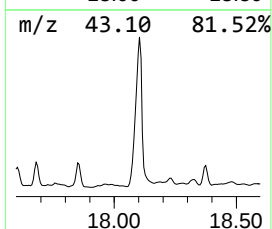
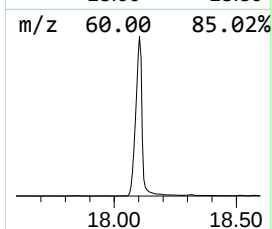
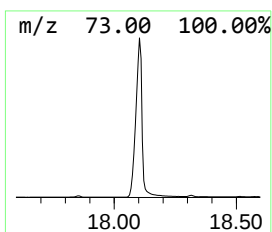
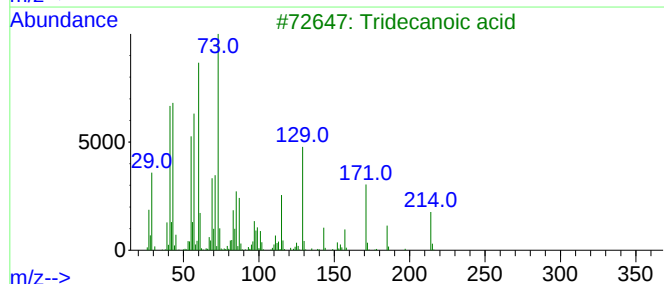
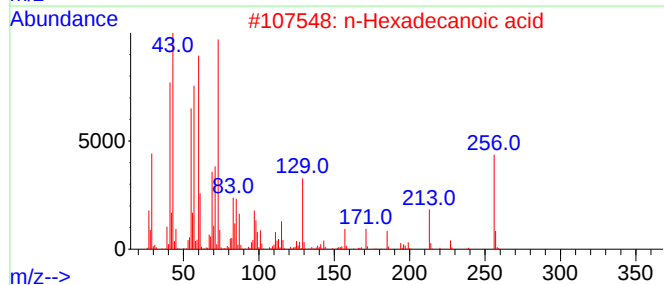
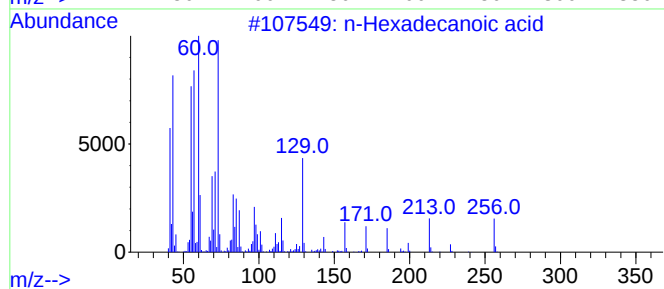
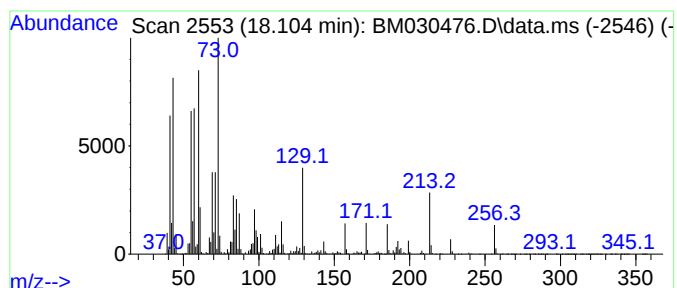
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.100	168.57 ng/ul	20987500	Phenanthrene-d10	17.186

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			Tridecanoic acid	214	C13H26O2	000638-53-9	92
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
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Instrument :
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ClientSampleId :
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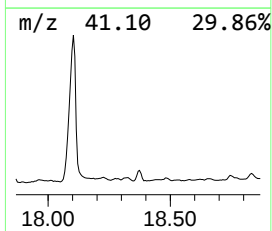
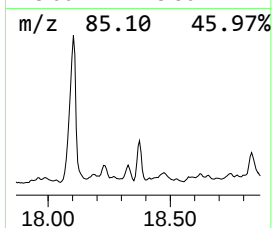
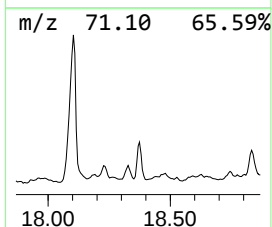
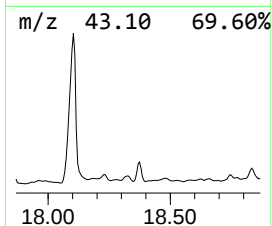
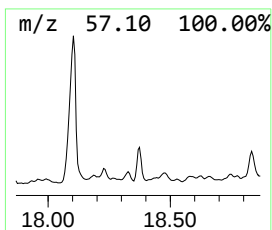
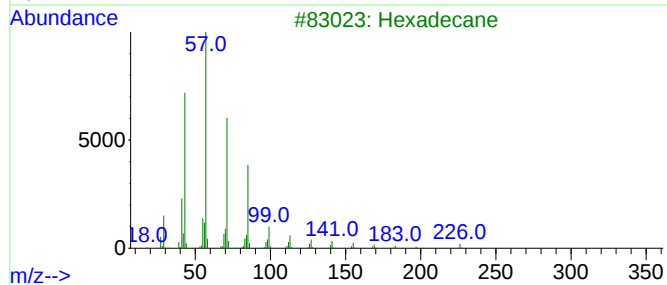
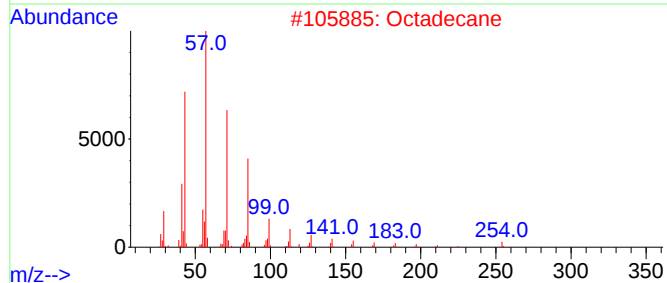
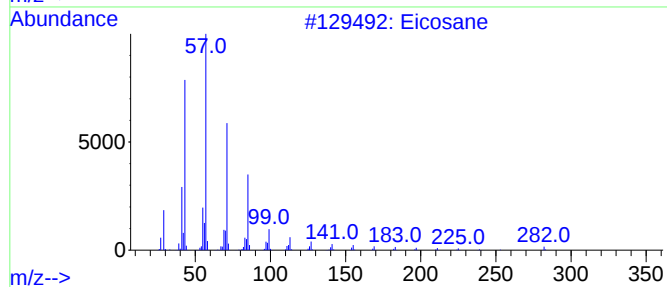
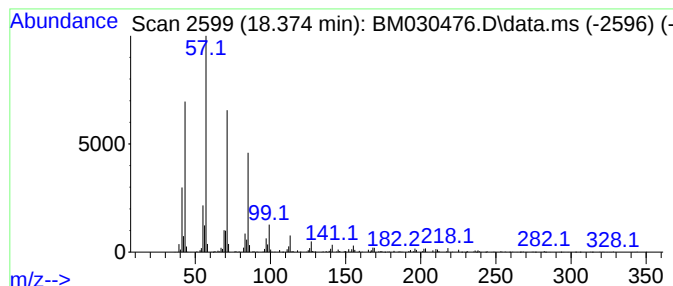
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.370	6.75 ng/ul	839977	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Octadecane	254	C18H38	000593-45-3	91
3			Hexadecane	226	C16H34	000544-76-3	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Sample : M2596-05
Misc :
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ClientSampleId :
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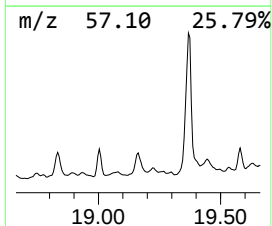
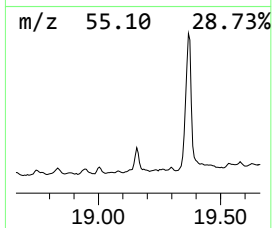
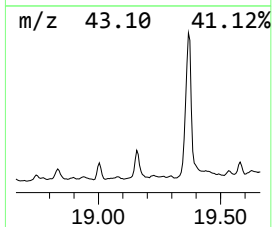
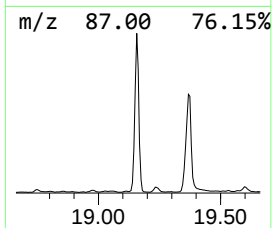
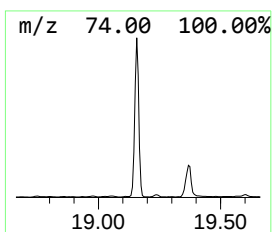
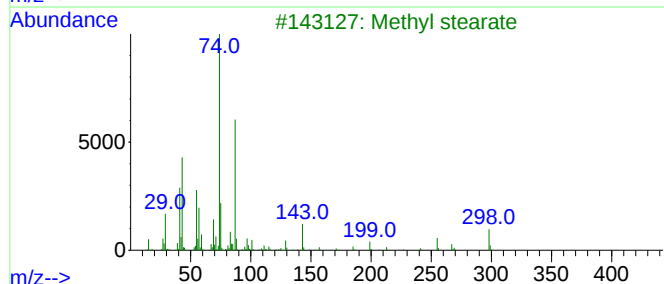
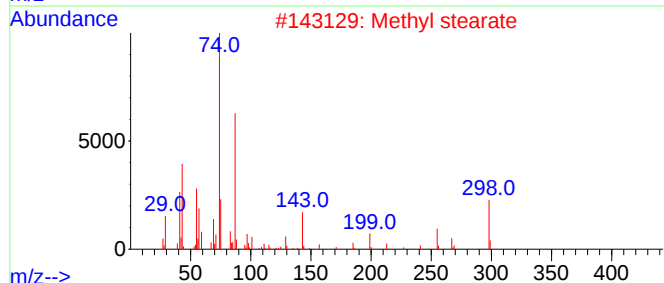
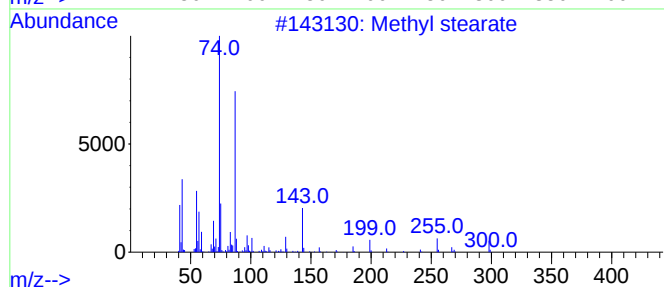
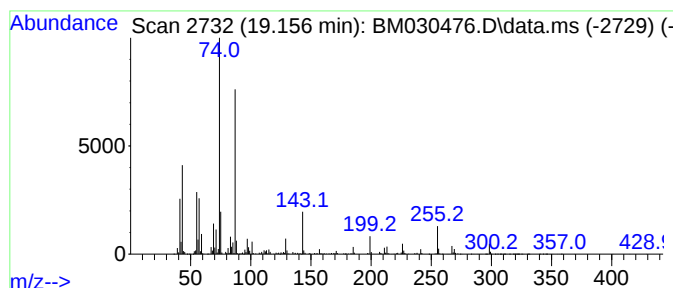
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 Methyl stearate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.160	23.99 ng/ul	2987430	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Methyl stearate	298	C19H38O2	000112-61-8	97
3			Methyl stearate	298	C19H38O2	000112-61-8	96
4			Methyl stearate	298	C19H38O2	000112-61-8	95
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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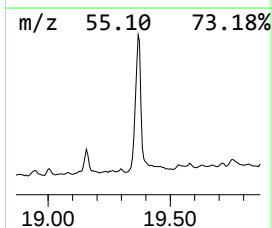
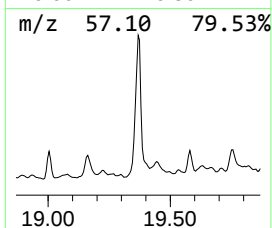
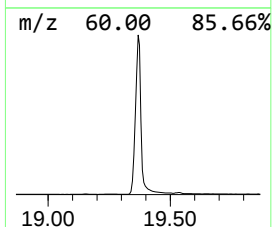
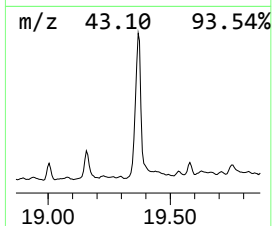
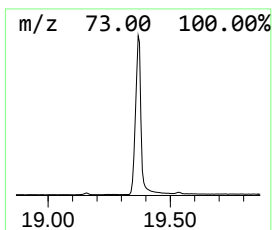
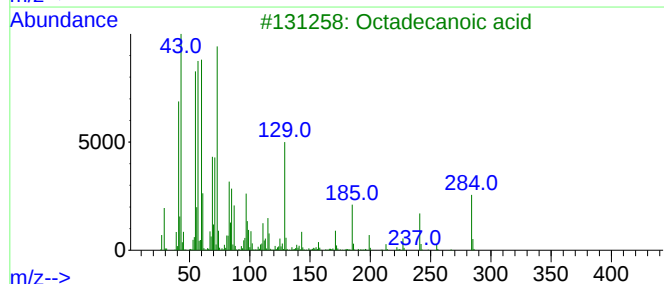
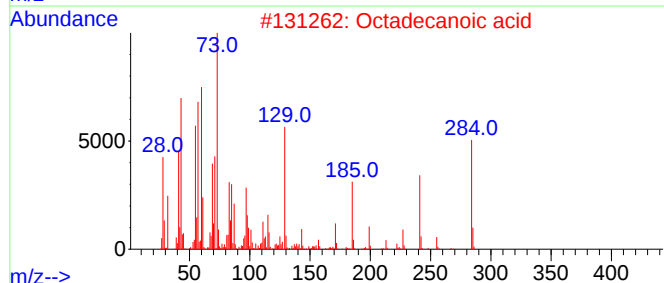
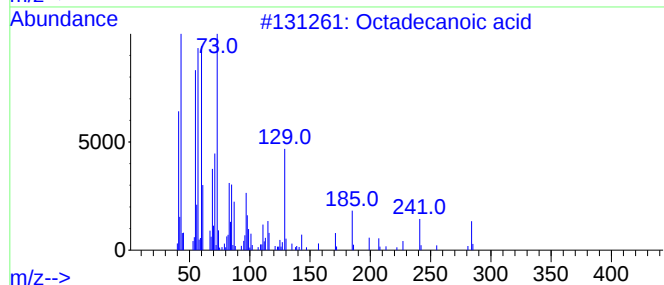
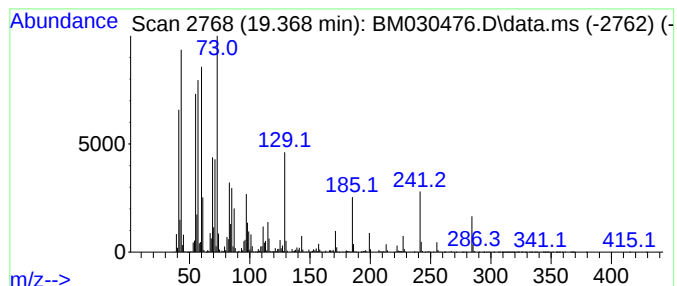
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.370	14.98 ng/ul	18818400	Chrysene-d12	21.350

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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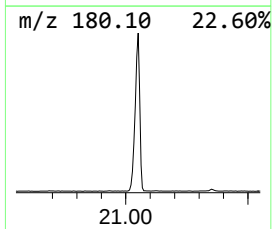
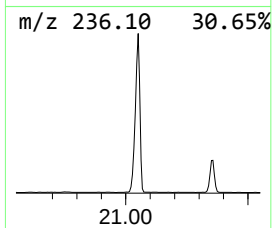
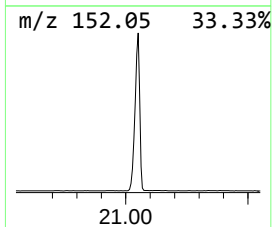
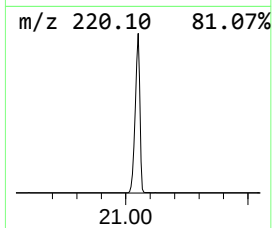
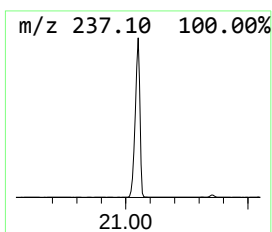
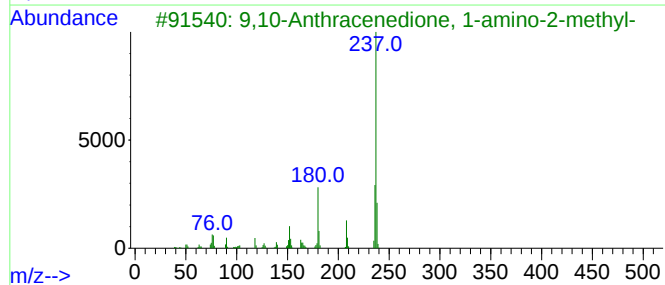
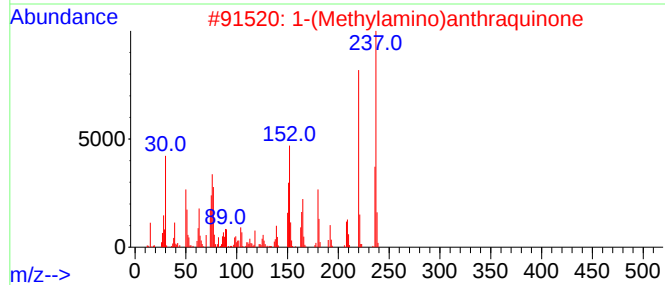
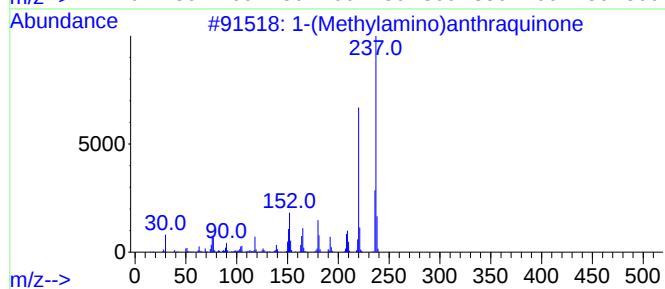
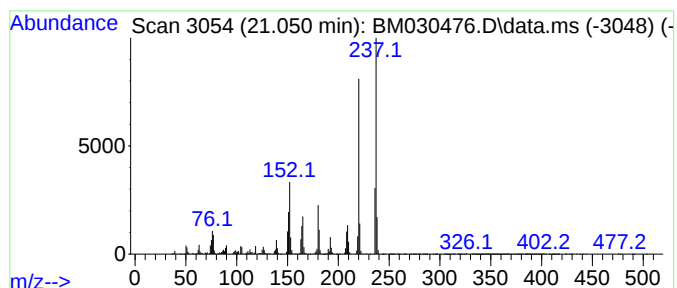
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 1-(Methylamino)anthraquinone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.050	20.00 ng/ul	25131900	Chrysene-d12	21.350

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(Methylamino)anthraquinone	237	C15H11NO2	000082-38-2	99		
2	1-(Methylamino)anthraquinone	237	C15H11NO2	000082-38-2	99		
3	9,10-Anthracenedione, 1-amino-2-...	237	C15H11NO2	000082-28-0	90		
4	7-Amino-3-phenylcoumarin	237	C15H11NO2	004108-61-6	38		
5	2-(p-(Dimethylamino)phenyl)benzi...	237	C15H15N3	002562-71-2	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030476.D
Acq On : 16 Jun 2021 21:10
Operator : CG/JU
Sample : M2596-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5Q1

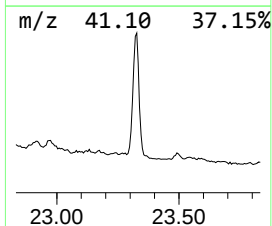
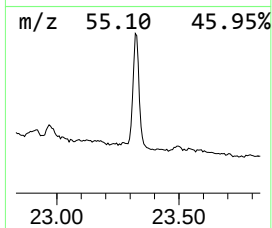
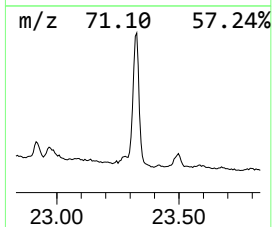
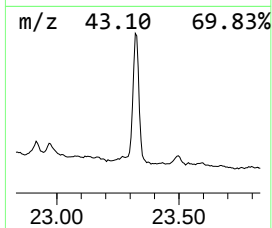
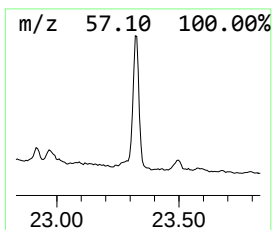
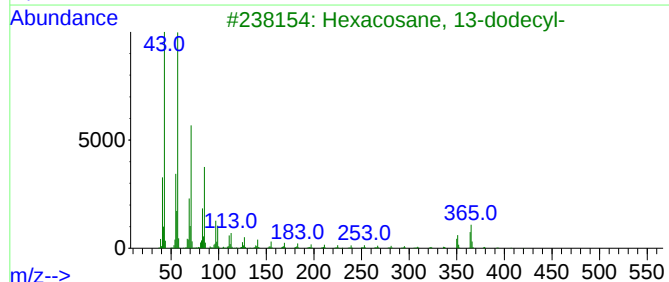
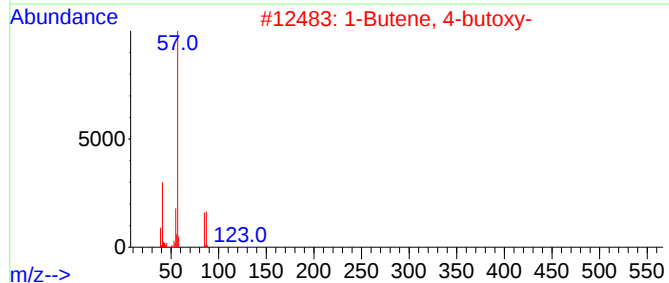
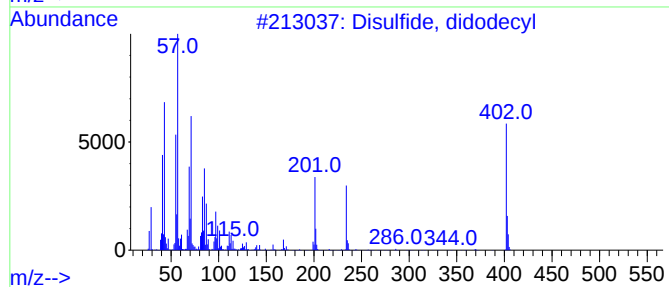
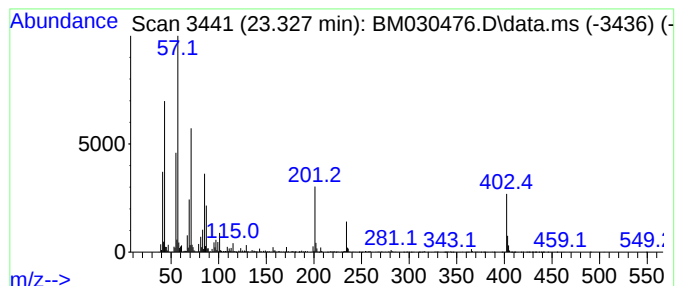
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 Disulfide, didodecyl Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.330	42.14 ng/ul	5345500	Perylene-d12	23.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Disulfide, didodecyl	402	C24H50S2	002757-37-1	78
2			1-Butene, 4-butoxy-	128	C8H16O	034061-76-2	22
3			Hexacosane, 13-dodecyl-	535	C38H78	055517-73-2	20
4			Threitol, 2-O-decyl-	262	C14H30O4	1000156-80-6	17
5			Hexacosane, 9-octyl-	479	C34H70	055429-83-9	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
 Data File : BM030476.D
 Acq On : 16 Jun 2021 21:10
 Operator : CG/JU
 Sample : M2596-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG5Q1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Propenoic aci...	3.970	16.6	ng/ul	816544	1	7.828	984094	20.0
unknown-01	8.920	3.4	ng/ul	168006	1	7.828	984094	20.0
(DEL) Alkane: S...	9.050	3.9	ng/ul	193189	1	7.828	984094	20.0
(DEL) Alkane: S...	9.960	2.6	ng/ul	236236	2	10.610	1827460	20.0
(DEL) Alkane: S...	10.040	5.1	ng/ul	463141	2	10.610	1827460	20.0
(DEL) Alkane: S...	10.150	2.6	ng/ul	234906	2	10.610	1827460	20.0
(DEL) Alkane: C...	10.460	12.8	ng/ul	1164690	2	10.610	1827460	20.0
unknown-02	10.830	2.7	ng/ul	247267	2	10.610	1827460	20.0
(DEL) Alkane: C...	11.100	4.5	ng/ul	413175	2	10.610	1827460	20.0
(DEL) Alkane: S...	11.270	2.1	ng/ul	193189	2	10.610	1827460	20.0
(DEL) Alkane: C...	11.310	9.5	ng/ul	864686	2	10.610	1827460	20.0
unknown-03	11.370	5.6	ng/ul	507843	2	10.610	1827460	20.0
Sulfurous acid,...	11.630	19.3	ng/ul	1760420	2	10.610	1827460	20.0
(DEL) Alkane: S...	11.900	5.6	ng/ul	513416	2	10.610	1827460	20.0
(DEL) Alkane: S...	12.160	2.7	ng/ul	245716	2	10.610	1827460	20.0
Hexanedioic aci...	12.340	8.3	ng/ul	755010	2	10.610	1827460	20.0
(DEL) Alkane: S...	12.670	5.8	ng/ul	597062	3	14.451	2076800	20.0
(DEL) Alkane: C...	12.730	7.6	ng/ul	784944	3	14.451	2076800	20.0
(DEL) Alkane: S...	12.770	2.9	ng/ul	297662	3	14.451	2076800	20.0
unknown-04	12.940	2.9	ng/ul	299535	3	14.451	2076800	20.0
(DEL) Alkane: S...	12.990	22.6	ng/ul	2346840	3	14.451	2076800	20.0
unknown-05	13.090	17.2	ng/ul	1782620	3	14.451	2076800	20.0
2-Decen-1-ol, (E)-	13.410	15.0	ng/ul	1560610	3	14.451	2076800	20.0
Naphthalene, 2,...	13.610	15.5	ng/ul	1610380	3	14.451	2076800	20.0
Naphthalene, 1,...	13.770	8.9	ng/ul	929513	3	14.451	2076800	20.0
unknown-06	13.970	3.3	ng/ul	342517	3	14.451	2076800	20.0
unknown-07	14.290	10.9	ng/ul	1136970	3	14.451	2076800	20.0
(DEL) Alkane: S...	14.350	12.1	ng/ul	1258100	3	14.451	2076800	20.0
Naphthalene, 2,...	14.850	24.4	ng/ul	2529410	3	14.451	2076800	20.0
Sulfurous acid,...	14.970	16.4	ng/ul	1701920	3	14.451	2076800	20.0
Naphthalene, 1,...	15.070	11.7	ng/ul	1214590	3	14.451	2076800	20.0
Sulfurous acid,...	15.300	8.1	ng/ul	843329	3	14.451	2076800	20.0
(DEL) Alkane: S...	15.700	40.1	ng/ul	4166100	3	14.451	2076800	20.0
unknown-08	15.860	5.7	ng/ul	706998	4	17.186	2490080	20.0
unknown-09	15.920	5.2	ng/ul	650910	4	17.186	2490080	20.0
(DEL) Alkane: S...	16.180	77.5	ng/ul	9646650	4	17.186	2490080	20.0
unknown-10	16.840	10.6	ng/ul	1314330	4	17.186	2490080	20.0
unknown-11	16.870	14.6	ng/ul	1818120	4	17.186	2490080	20.0
(DEL) Alkane: S...	17.000	35.5	ng/ul	4421380	4	17.186	2490080	20.0
unknown-12	17.360	5.7	ng/ul	705924	4	17.186	2490080	20.0
unknown-13	17.420	8.6	ng/ul	1068790	4	17.186	2490080	20.0
Hexadecane, 1-i...	17.600	15.4	ng/ul	1916070	4	17.186	2490080	20.0
(DEL) Alkane: S...	17.680	11.1	ng/ul	1375490	4	17.186	2490080	20.0
Pentadecanoic a...	17.850	21.6	ng/ul	2693430	4	17.186	2490080	20.0
n-Hexadecanoic ...	18.100	168.6	ng/ul	20987500	4	17.186	2490080	20.0
(DEL) Alkane: S...	18.370	6.8	ng/ul	839977	4	17.186	2490080	20.0
Methyl stearate	19.160	24.0	ng/ul	2987430	4	17.186	2490080	20.0
Octadecanoic acid	19.370	15.0	ng/ul	18818400	5	21.350	25131900	20.0
1-(Methylamino)...	21.050	20.0	ng/ul	25131900	5	21.350	25131900	20.0
Disulfide, dido...	23.330	42.1	ng/ul	5345500	6	23.627	2536940	20.0