

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
 Data File : BM030477.D
 Acq On : 16 Jun 2021 21:46
 Operator : CG/JU
 Sample : M2596-05DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG5Q1DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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: BM030477.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.322	35	40	75	rBV	676335	1498047	31.79%	3.391%
2	3.816	121	124	131	rVB	18143	24366	0.52%	0.055%
3	3.910	133	140	144	rBV	69698	125352	2.66%	0.284%
4	4.040	158	162	169	rVB	16605	25703	0.55%	0.058%
5	4.587	251	255	262	rVB2	14151	22324	0.47%	0.051%
6	5.481	402	407	416	rVB	16363	34333	0.73%	0.078%
7	6.998	656	665	678	rBV	81810	161145	3.42%	0.365%
8	7.157	685	692	701	rBV	62979	108680	2.31%	0.246%
9	7.310	713	718	721	rBV2	18069	25742	0.55%	0.058%
10	7.363	721	727	734	rVB	87248	150646	3.20%	0.341%
11	7.822	798	805	814	rBV	591396	983251	20.86%	2.226%
12	8.528	920	925	932	rVV	92920	161820	3.43%	0.366%
13	8.916	985	991	996	rBV7	10740	23540	0.50%	0.053%
14	8.981	996	1002	1008	rVV	72329	128641	2.73%	0.291%
15	9.045	1008	1013	1018	rVB2	19884	34509	0.73%	0.078%
16	9.298	1048	1056	1062	rVB2	10059	27611	0.59%	0.063%
17	9.463	1078	1084	1088	rVV4	14490	25390	0.54%	0.057%
18	9.698	1116	1124	1129	rBV	65832	119213	2.53%	0.270%
19	9.745	1129	1132	1136	rVV5	14633	21951	0.47%	0.050%
20	9.798	1136	1141	1148	rVB8	18997	35131	0.75%	0.080%
21	9.892	1148	1157	1161	rBV2	25907	63474	1.35%	0.144%
22	9.963	1161	1169	1174	rBV3	27097	49682	1.05%	0.112%
23	10.039	1174	1182	1193	rBV3	33540	95434	2.03%	0.216%
24	10.139	1195	1199	1204	rBV	30848	45171	0.96%	0.102%
25	10.239	1211	1216	1225	rVB	91919	173227	3.68%	0.392%
26	10.392	1238	1242	1247	rBV8	15243	28176	0.60%	0.064%
27	10.463	1248	1254	1259	rBV	142058	219160	4.65%	0.496%
28	10.610	1268	1279	1283	rBV	778978	1489277	31.60%	3.371%
29	10.651	1283	1286	1296	rVB	1017315	1565847	33.23%	3.545%
30	10.769	1296	1306	1313	rVB3	137139	309054	6.56%	0.700%
31	10.833	1313	1317	1326	rBV8	23509	51131	1.09%	0.116%
32	10.904	1326	1329	1335	rVB7	19054	27050	0.57%	0.061%
33	11.022	1346	1349	1354	rVB7	16032	26483	0.56%	0.060%
34	11.098	1356	1362	1369	rVB3	38417	79047	1.68%	0.179%
35	11.186	1372	1377	1381	rVB7	23612	38836	0.82%	0.088%
36	11.280	1388	1393	1394	rBV4	30391	47763	1.01%	0.108%

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Integrator: RTE

Smoother: OFF

Filtering: 5

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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
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37	11.310	1394	1398	1405	rVB7	42191	82158	1.74%	0.186%
38	11.375	1405	1409	1415	rVB5	38725	72497	1.54%	0.164%
39	11.445	1416	1421	1428	rVB5	54245	92649	1.97%	0.210%
40	11.527	1429	1435	1442	rVB8	42439	99476	2.11%	0.225%

41	11.633	1447	1453	1457	rBV	191125	319943	6.79%	0.724%
42	11.804	1477	1482	1489	rVB4	52113	97429	2.07%	0.221%
43	11.898	1491	1498	1502	rBV7	41939	81311	1.73%	0.184%
44	11.939	1502	1505	1509	rVV3	16720	21917	0.47%	0.050%
45	12.163	1540	1543	1547	rVB4	39388	54580	1.16%	0.124%

46	12.251	1552	1558	1567	rVB5	79188	209740	4.45%	0.475%
47	12.339	1569	1573	1580	rBV6	65673	107211	2.28%	0.243%
48	12.480	1592	1597	1600	rBV4	50350	91993	1.95%	0.208%
49	12.522	1601	1604	1609	rVB4	72267	113258	2.40%	0.256%
50	12.663	1623	1628	1632	rBV4	69043	115755	2.46%	0.262%

51	12.727	1632	1639	1644	rVV6	55187	139489	2.96%	0.316%
52	12.774	1644	1647	1650	rBV2	45013	57504	1.22%	0.130%
53	12.939	1671	1675	1678	rBV6	43919	86021	1.83%	0.195%
54	12.986	1679	1683	1689	rVB	286531	422162	8.96%	0.956%
55	13.057	1689	1695	1699	rBV	167016	270790	5.75%	0.613%

56	13.233	1721	1725	1728	rBV4	89760	150076	3.18%	0.340%
57	13.410	1744	1755	1758	rBV5	167459	394426	8.37%	0.893%
58	13.610	1783	1789	1796	rBV	110781	297407	6.31%	0.673%
59	13.769	1813	1816	1821	rVB	106727	154022	3.27%	0.349%
60	13.821	1822	1825	1827	rVV2	61323	74770	1.59%	0.169%

61	13.857	1828	1831	1836	rVV2	162698	289691	6.15%	0.656%
62	13.904	1836	1839	1841	rVV	244549	315136	6.69%	0.713%
63	13.933	1841	1844	1848	rVB	398828	490770	10.41%	1.111%
64	13.974	1848	1851	1854	rBV4	57156	64191	1.36%	0.145%
65	14.098	1868	1872	1875	rBV5	66667	112531	2.39%	0.255%

66	14.145	1875	1880	1884	rVV	223212	328758	6.98%	0.744%
67	14.286	1900	1904	1909	rVB3	120586	205089	4.35%	0.464%
68	14.351	1910	1915	1920	rBV4	123705	220225	4.67%	0.499%
69	14.451	1923	1932	1938	rVV	1212122	1988047	42.19%	4.500%
70	14.663	1964	1968	1976	rVB4	135465	341991	7.26%	0.774%

71	14.845	1995	1999	2006	rVB3	239382	447545	9.50%	1.013%
72	14.921	2010	2012	2016	rVB	121159	139636	2.96%	0.316%
73	14.974	2016	2021	2031	rVB6	151949	314117	6.67%	0.711%
74	15.068	2033	2037	2040	rBV	112142	188063	3.99%	0.426%
75	15.298	2073	2076	2080	rVB	114421	143853	3.05%	0.326%

76	15.398	2090	2093	2096	rBV4	56196	84279	1.79%	0.191%
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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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 Title : SVOA CALIBRATION

77	15.439	2096	2100	2104	rVB	276982	364588	7.74%	0.825%
78	15.663	2134	2138	2140	rBV4	73291	121143	2.57%	0.274%
79	15.704	2140	2145	2151	rVB2	498146	766841	16.27%	1.736%
80	15.857	2168	2171	2175	rBV4	71024	116519	2.47%	0.264%
81	15.910	2178	2180	2184	rVB2	90783	112015	2.38%	0.254%
82	16.180	2218	2226	2238	rVB	954789	1671765	35.48%	3.784%
83	16.839	2335	2338	2340	rBV2	179498	231880	4.92%	0.525%
84	16.862	2340	2342	2347	rVB2	269954	302688	6.42%	0.685%
85	16.945	2353	2356	2359	rVV	134461	161839	3.43%	0.366%
86	16.998	2361	2365	2376	rVB	633082	1084389	23.01%	2.455%
87	17.186	2392	2397	2402	rBV	1477956	2131681	45.24%	4.825%
88	17.280	2410	2413	2419	rVB	333882	465283	9.87%	1.053%
89	17.598	2463	2467	2477	rBV2	198516	366300	7.77%	0.829%
90	17.680	2478	2481	2487	rVB2	192494	232556	4.93%	0.526%
91	17.851	2506	2510	2515	rVB	368193	476274	10.11%	1.078%
92	18.080	2543	2549	2561	rBV2	2081581	3127823	66.37%	7.080%
93	18.368	2596	2598	2604	rVB	126936	154079	3.27%	0.349%
94	19.156	2728	2732	2740	rVB2	406112	542364	11.51%	1.228%
95	19.351	2761	2765	2774	rBV	1611077	2389377	50.70%	5.409%
96	19.568	2798	2802	2815	rBV4	420406	936861	19.88%	2.121%
97	21.033	3047	3051	3059	rBV	3603477	4712457	100.00%	10.667%
98	21.350	3100	3105	3109	rBV	2048782	2691754	57.12%	6.093%
99	23.321	3437	3440	3452	rVB	543802	903139	19.16%	2.044%
100	23.621	3486	3491	3502	rVB	1489345	2785945	59.12%	6.306%

Sum of corrected areas: 44176273

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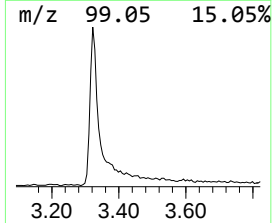
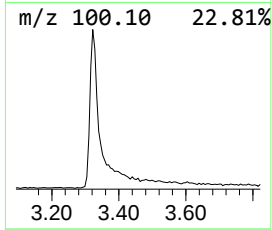
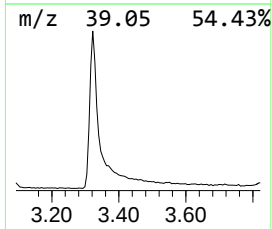
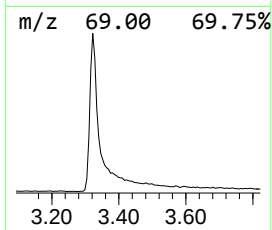
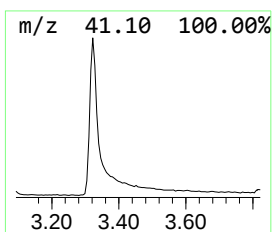
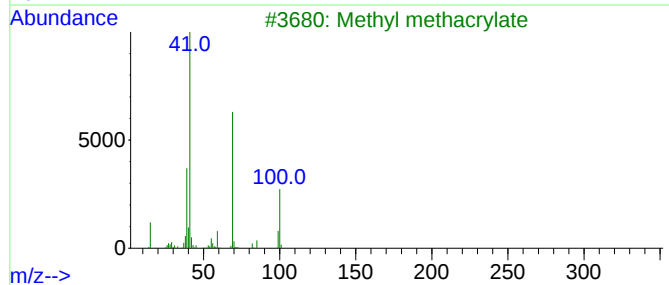
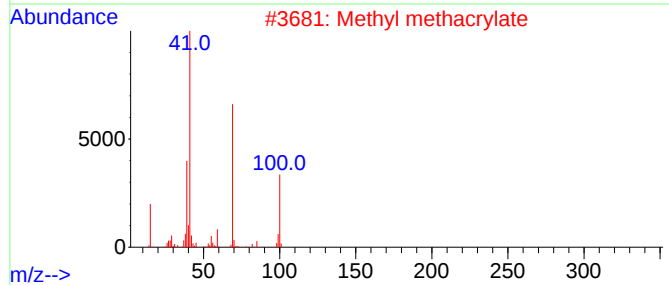
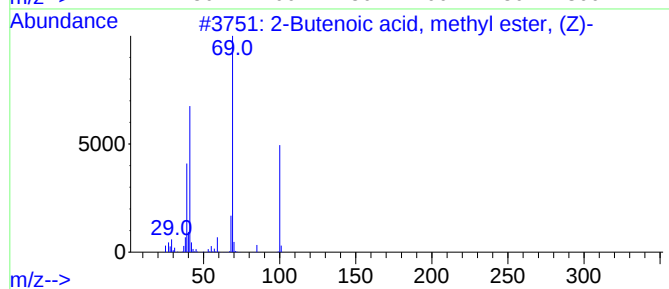
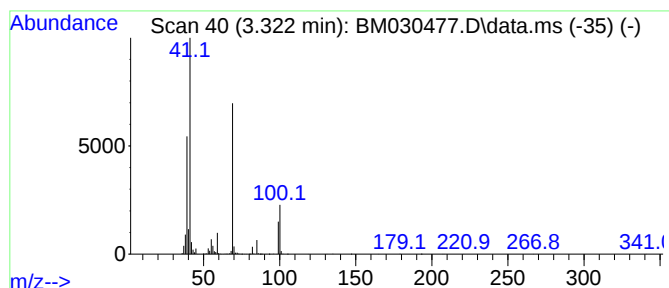
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Butenoic acid, methyl est... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.320	30.47 ng/ul	1498050	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butenoic acid, methyl ester, (Z)-	100	C5H8O2	004358-59-2	86
2			Methyl methacrylate	100	C5H8O2	000080-62-6	80
3			Methyl methacrylate	100	C5H8O2	000080-62-6	64
4			Methyl methacrylate	100	C5H8O2	000080-62-6	60
5			2-Propenoic acid, 2-methyl-, 2-h...	144	C7H12O3	000923-26-2	56



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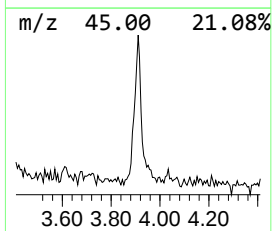
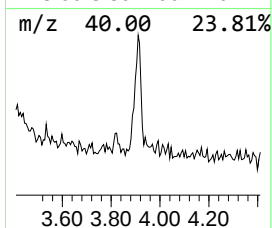
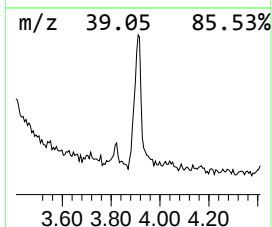
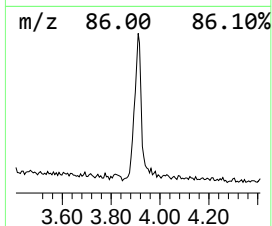
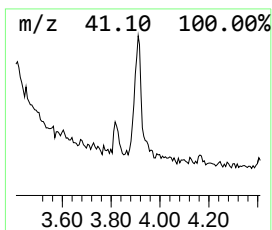
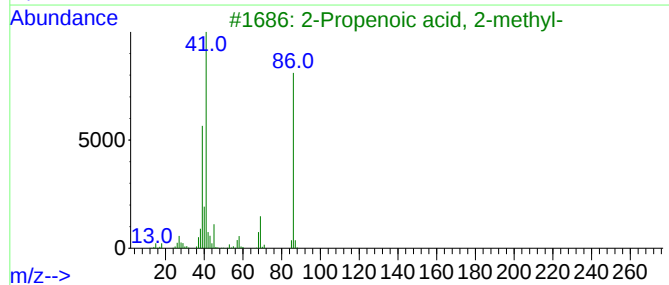
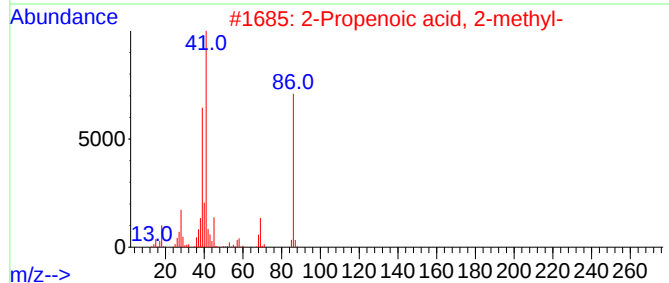
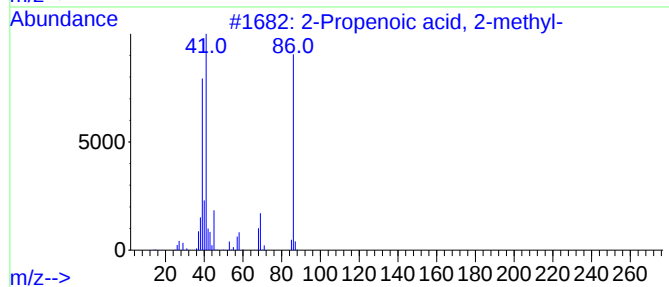
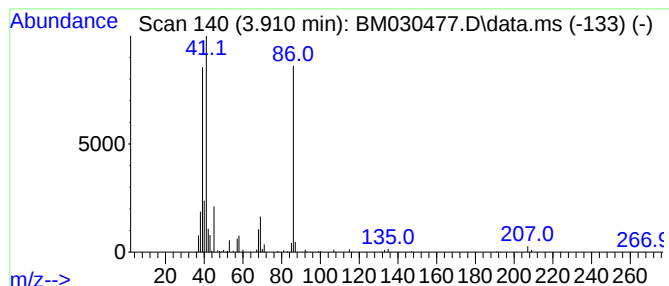
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propenoic acid, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.910	2.55 ng/ul	125352	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	91
3			2-Propenoic acid, 2-methyl-	86	C4H6O2	000079-41-4	91
4			Crotonic acid	86	C4H6O2	003724-65-0	90
5			Crotonic acid	86	C4H6O2	003724-65-0	90



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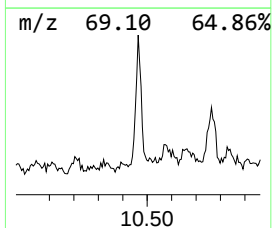
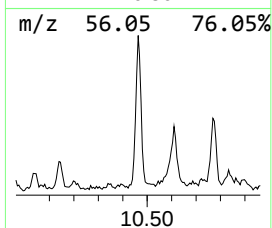
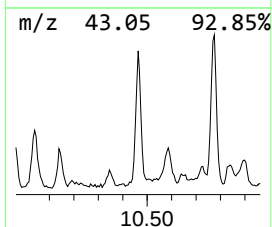
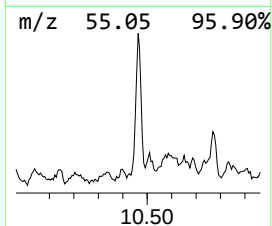
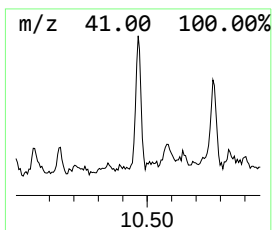
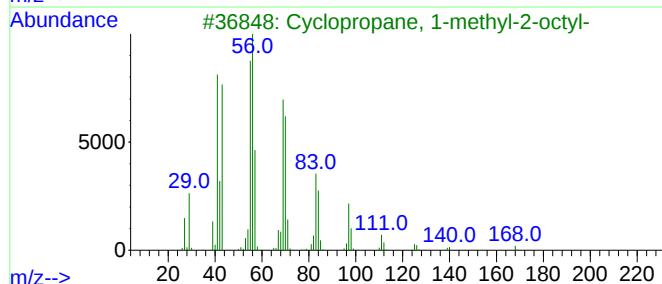
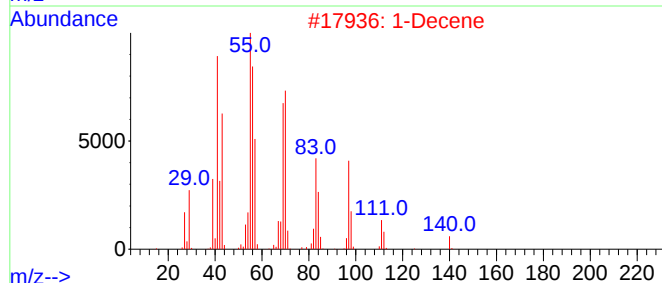
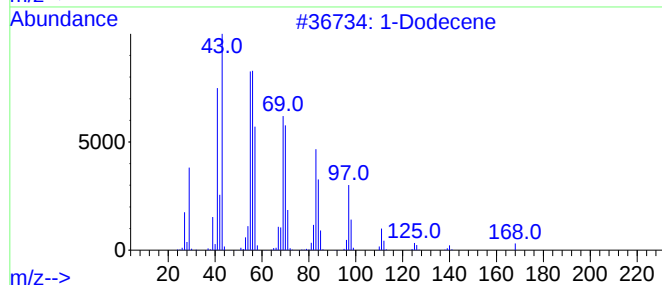
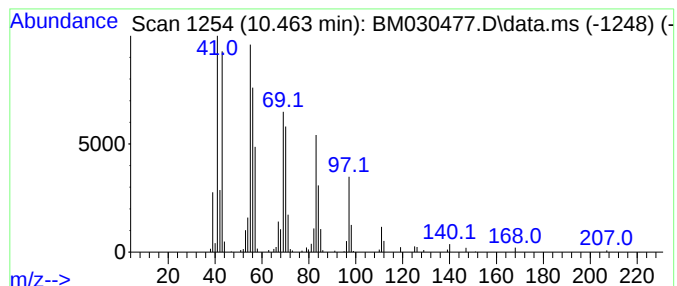
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.460	2.94 ng/ul	219160	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Dodecene	168	C12H24	000112-41-4	97
2			1-Decene	140	C10H20	000872-05-9	93
3			Cyclopropane, 1-methyl-2-octyl-	168	C12H24	037617-26-8	93
4			1-Tridecene	182	C13H26	002437-56-1	91
5			E-11,13-Tetradecadien-1-ol	210	C14H26O	1000131-00-3	91



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BNA_M
ClientSampleId :
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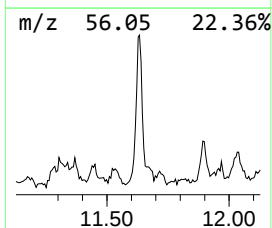
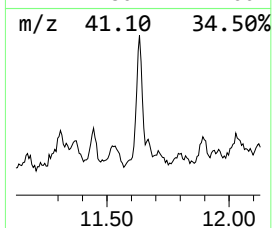
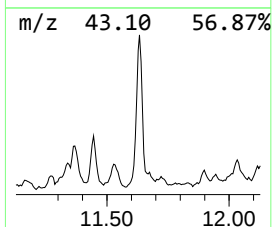
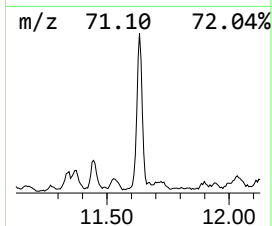
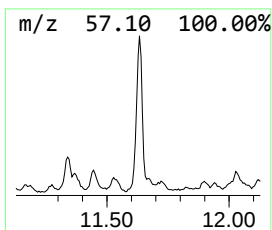
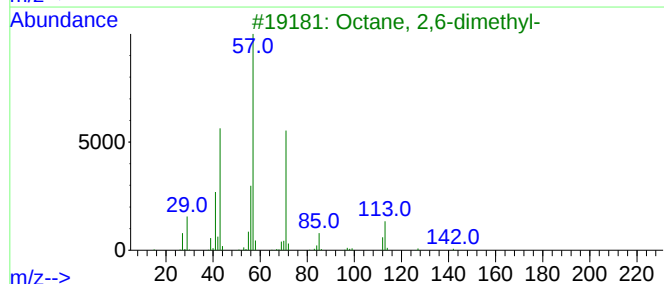
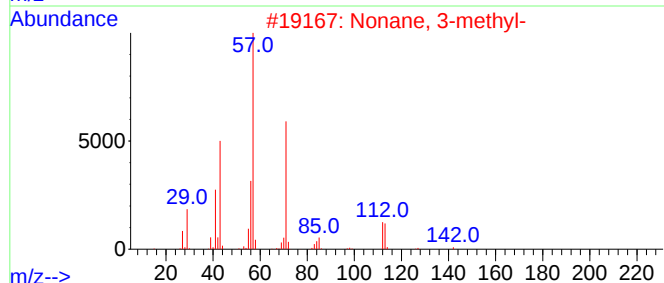
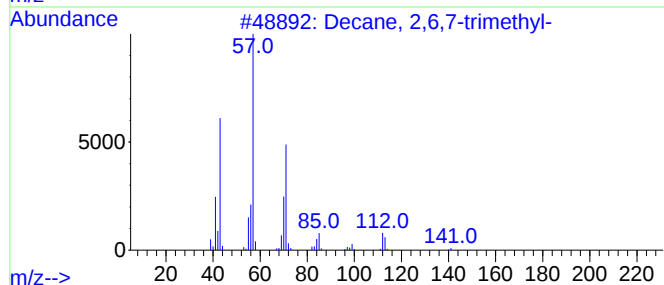
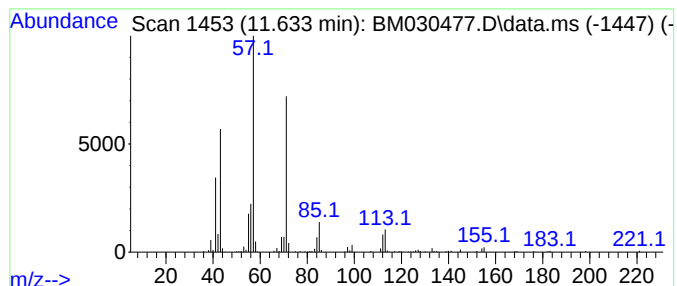
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.630	4.30 ng/ul	319943	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	90
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	78
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	78
4		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64
5		Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	64



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Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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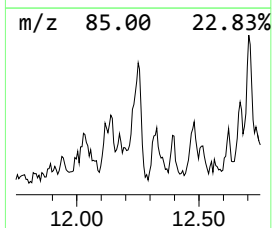
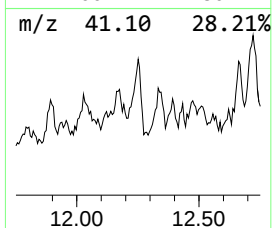
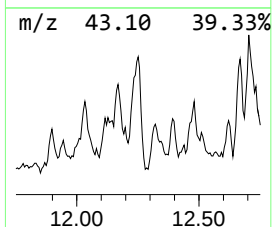
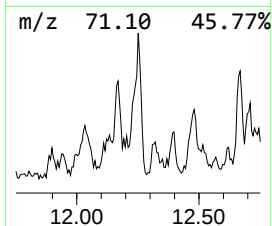
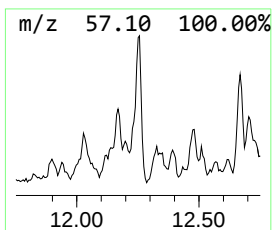
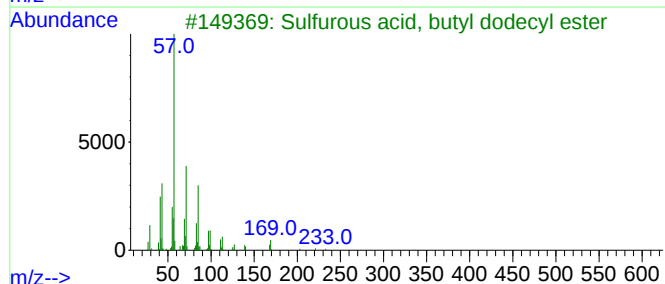
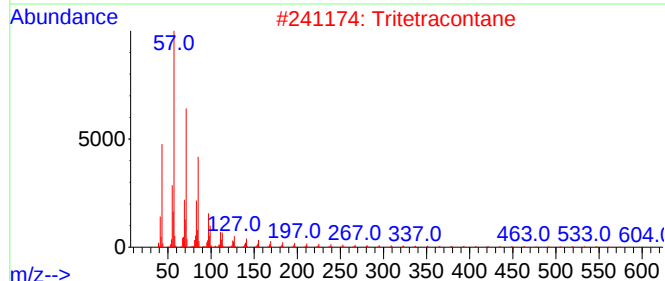
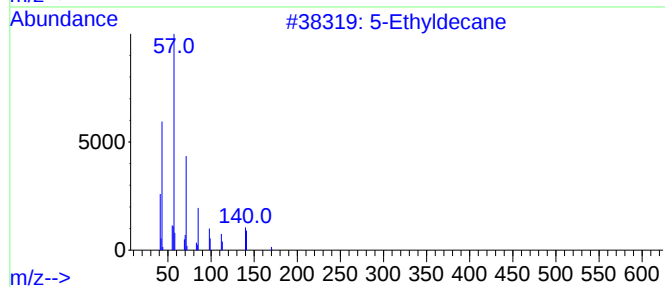
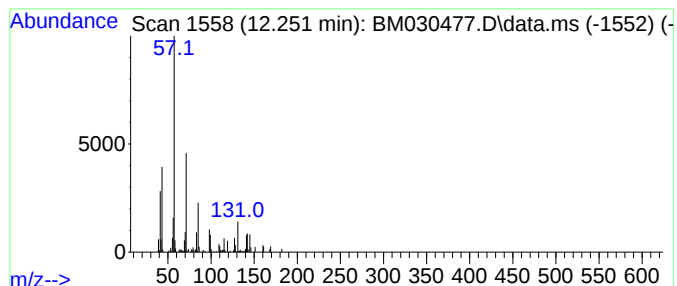
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	2.82 ng/ul	209740	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Ethyldecane	170	C12H26	017302-36-2	59
2			Tritetracontane	605	C43H88	007098-21-7	50
3			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	50
4			Decane, 1-bromo-2-methyl-	234	C11H23Br	127839-47-8	47
5			Dodecane, 6-methyl-	184	C13H28	006044-71-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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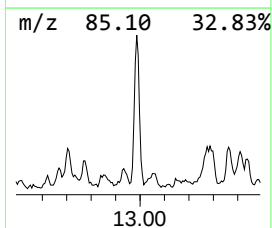
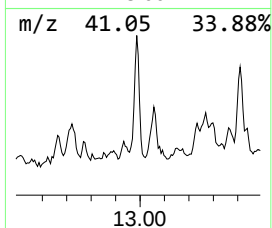
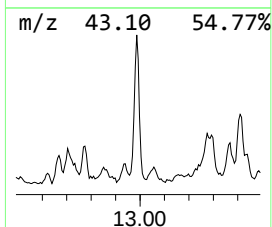
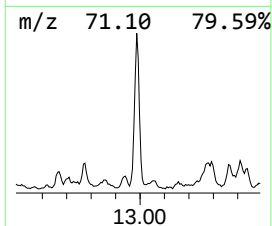
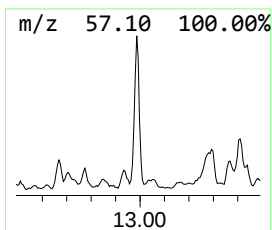
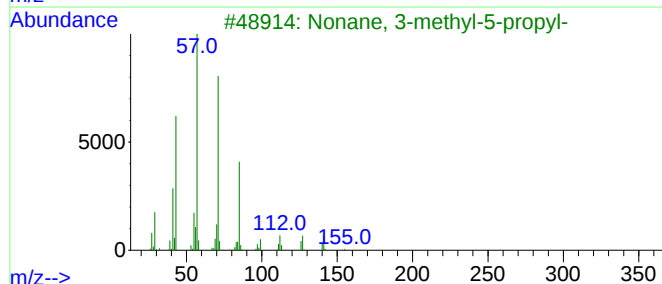
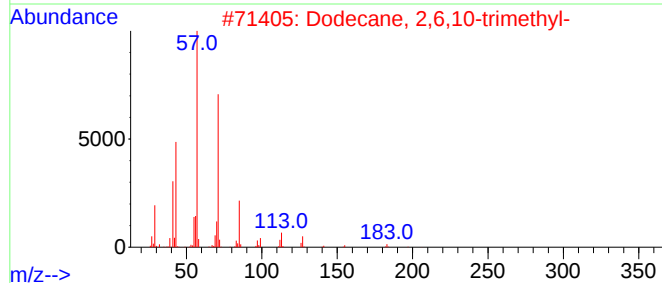
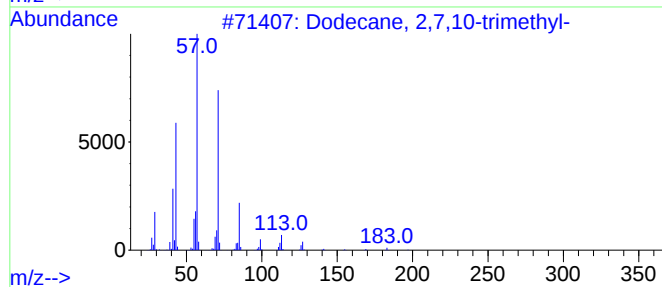
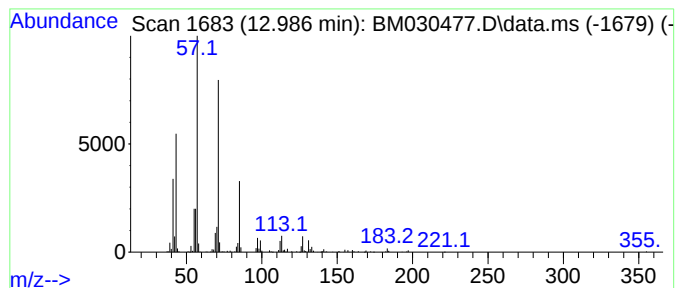
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.990	4.25 ng/ul	422162	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			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
4			Bacchotricuneatin c	342	C20H22O5	066563-30-2	72
5			Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
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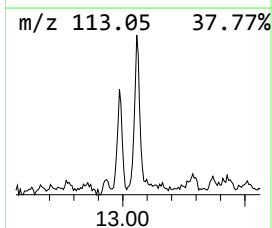
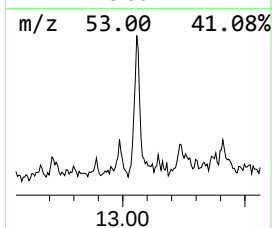
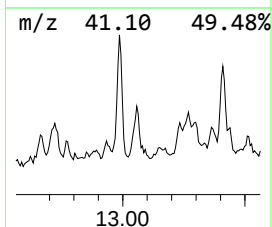
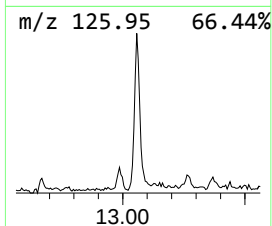
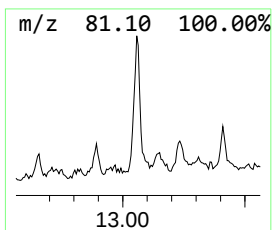
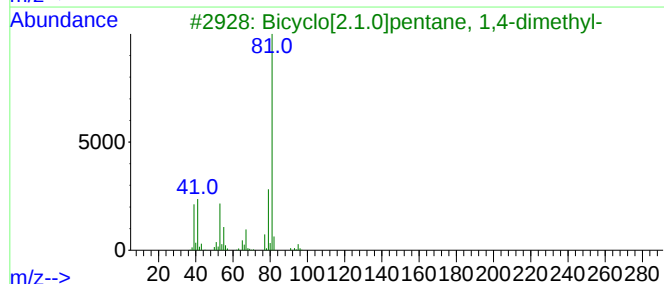
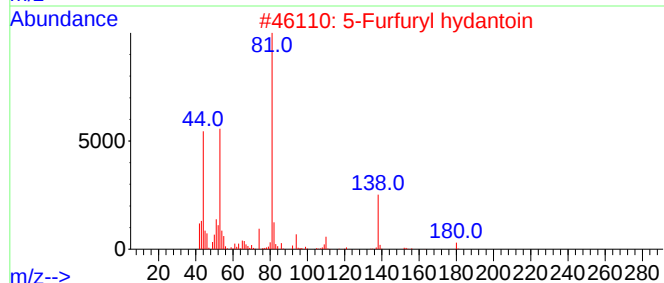
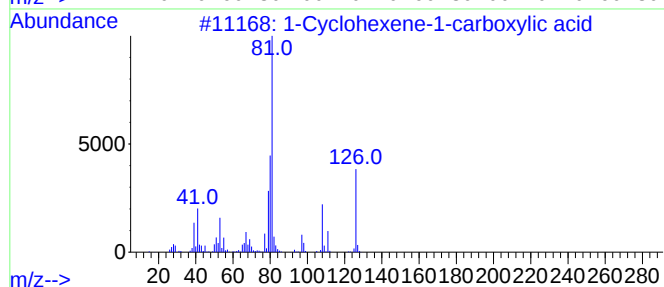
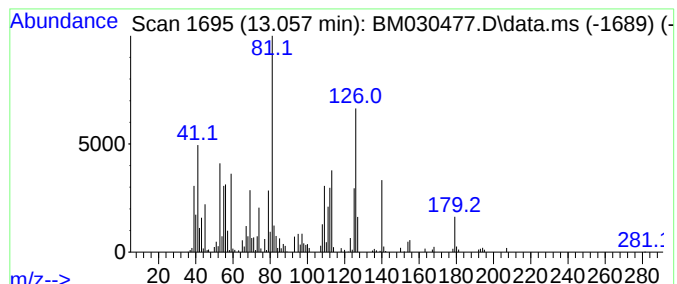
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Cyclohexene-1-carboxylic ... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.060	2.72 ng/ul	270790	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Cyclohexene-1-carboxylic acid	126	C7H10O2	000636-82-8	58
2			5-Furfuryl hydantoin	180	C8H8N2O3	004349-14-8	43
3			Bicyclo[2.1.0]pentane, 1,4-dimet...	96	C7H12	017065-18-8	43
4			2,4-Pentadienoic acid, 1-cyclope...	178	C10H10O3	1000150-80-4	43
5			1-Cyclohexene-1-carboxylic acid	126	C7H10O2	000636-82-8	42



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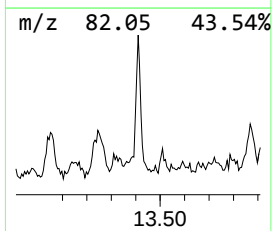
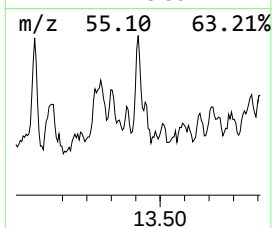
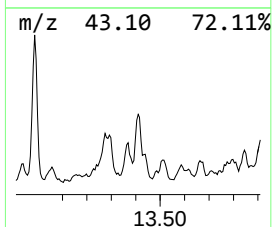
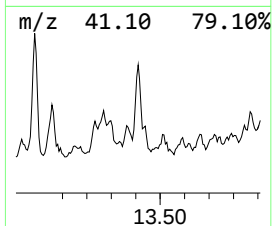
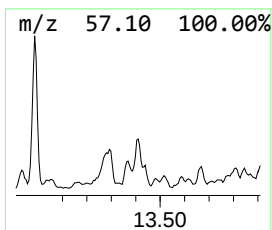
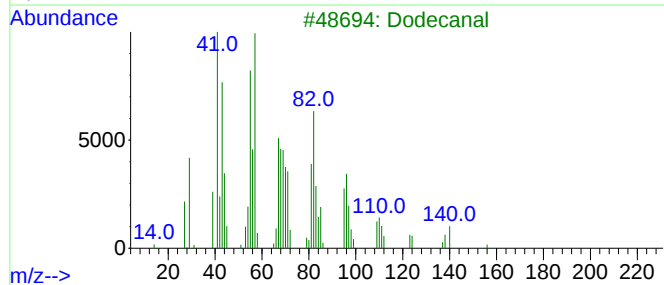
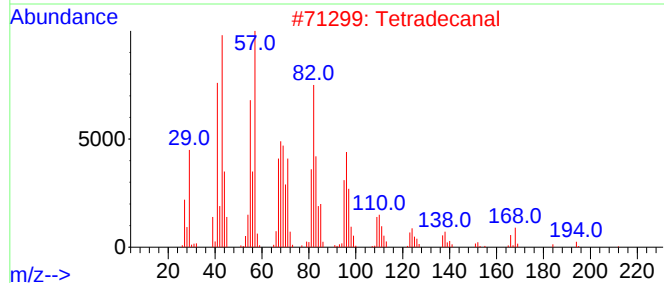
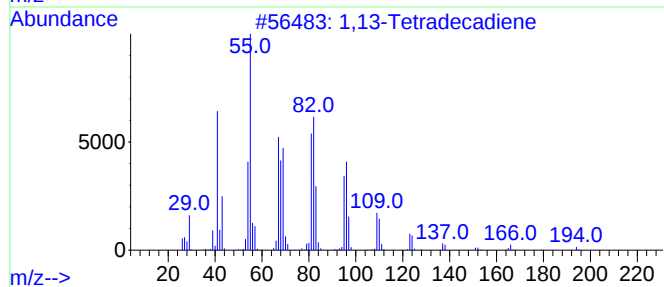
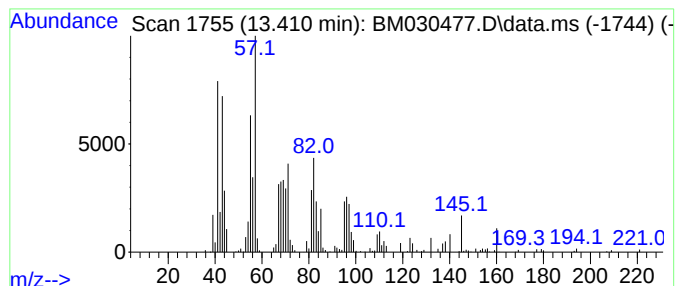
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,13-Tetradecadiene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.410	3.97 ng/ul	394426	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,13-Tetradecadiene	194	C14H26	021964-49-8	58
2			Tetradecanal	212	C14H28O	000124-25-4	58
3			Dodecanal	184	C12H24O	000112-54-9	58
4			Heptadecanal	254	C17H34O	1000376-70-0	58
5			2-Decen-1-ol, (E)-	156	C10H20O	018409-18-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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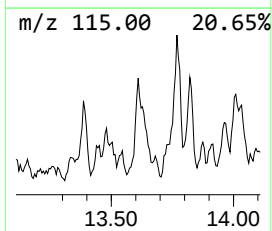
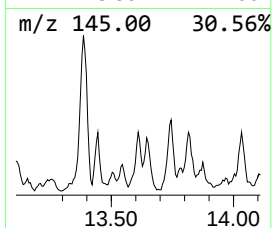
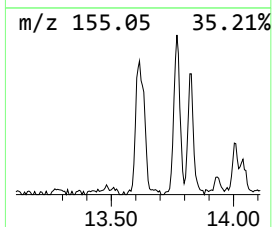
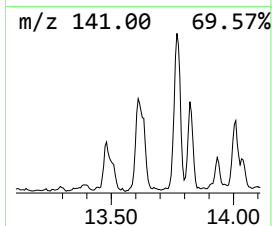
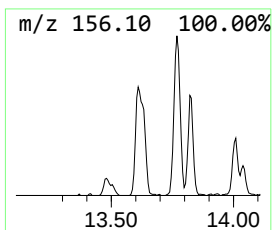
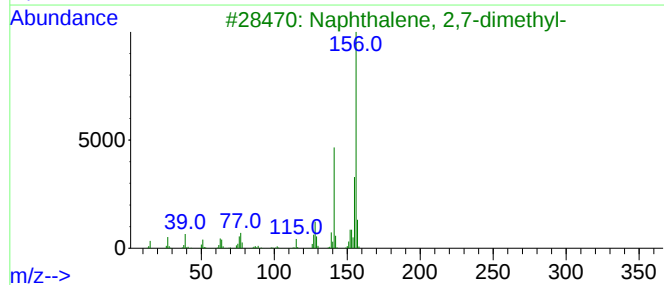
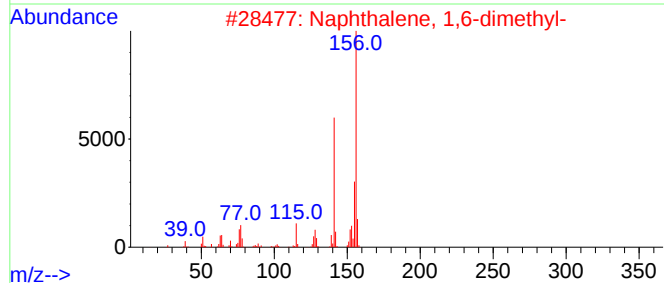
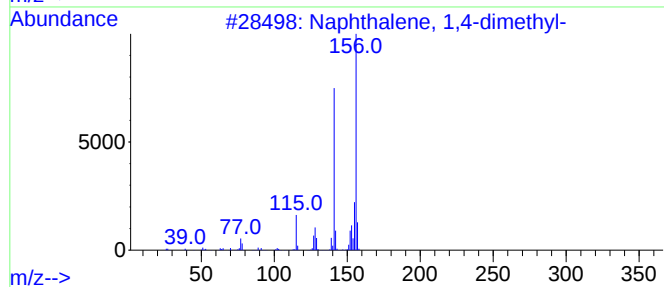
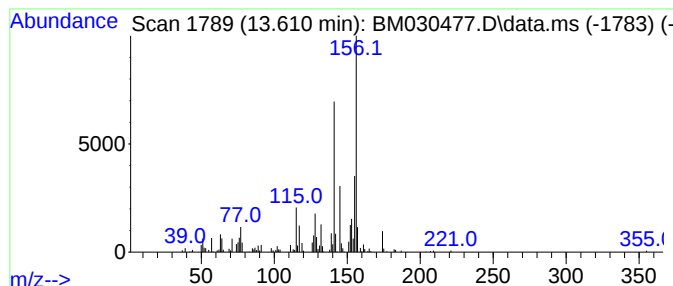
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalene, 1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.610	2.99 ng/ul	297407	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94



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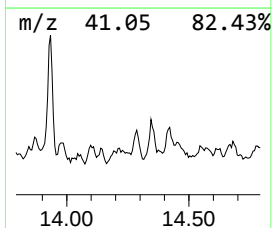
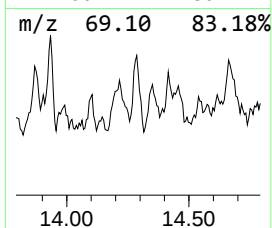
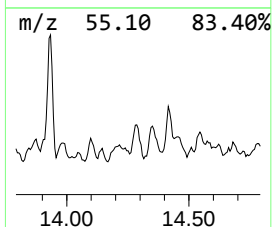
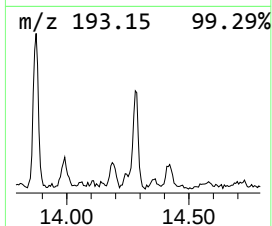
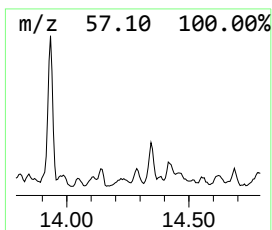
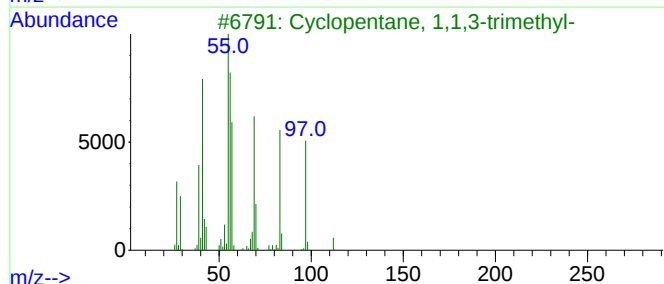
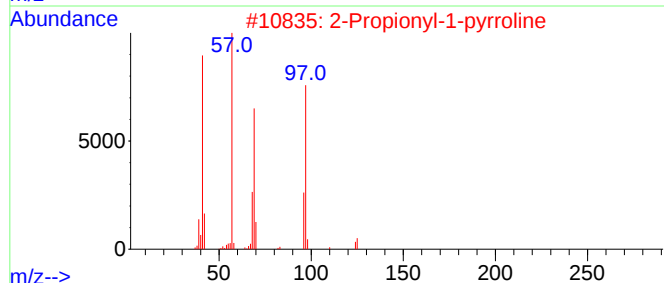
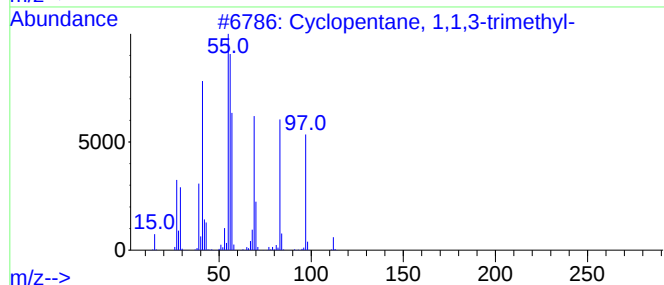
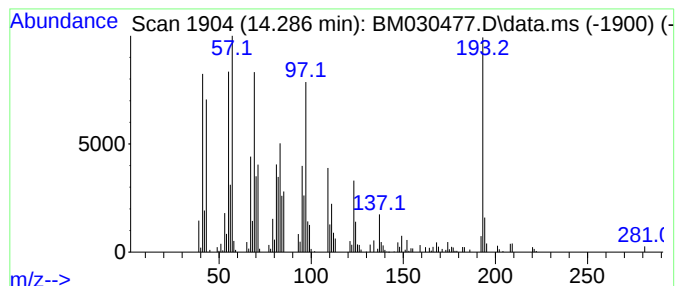
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic14.29 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.290	2.06 ng/ul	205089	Acenaphthene-d10			14.451
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1,3-trimethyl-		112	C8H16	004516-69-2	30
2	2-Propionyl-1-pyrroline		125	C7H11NO	1000298-55-4	30
3	Cyclopentane, 1,1,3-trimethyl-		112	C8H16	004516-69-2	30
4	4H-Pyran-4-one, 2,6-dimethyl-		124	C7H8O2	001004-36-0	25
5	Thieno[2,3-d]-1,3-thiaselenol-2-...		238	C5H2S3Se	081803-09-0	25



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Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

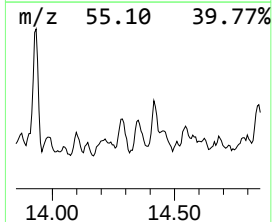
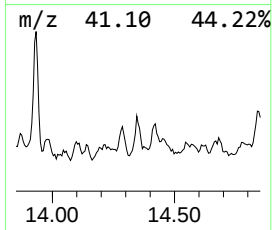
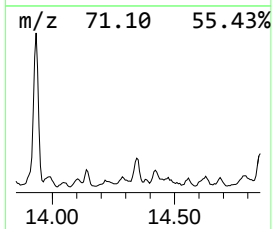
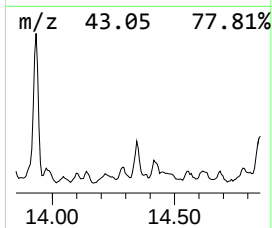
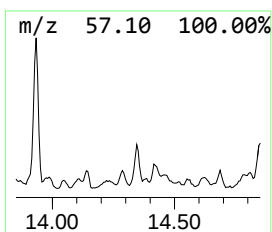
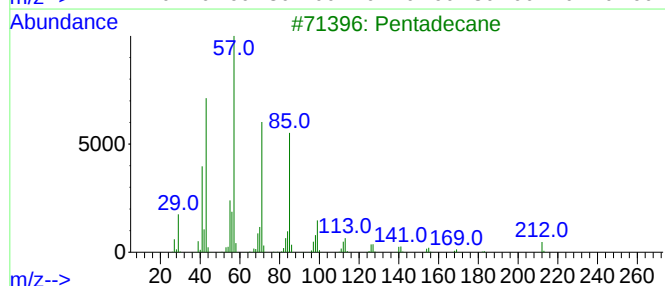
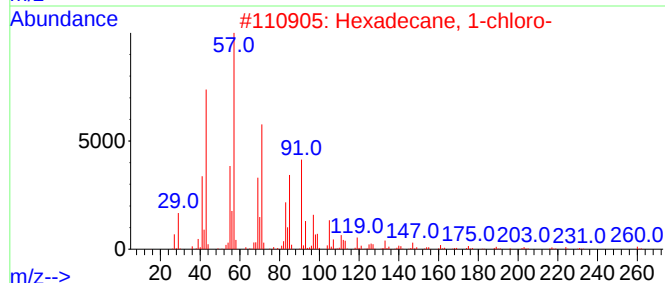
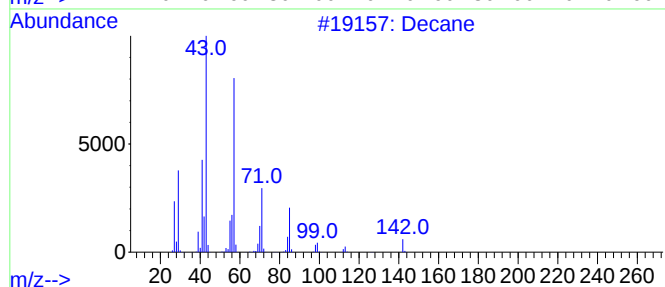
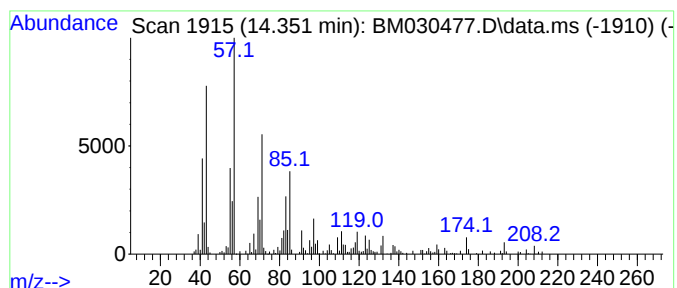
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.350	2.22 ng/ul	220225	Acenaphthene-d10		14.451	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	92
2	Hexadecane, 1-chloro-		260	C16H33Cl	004860-03-1	80
3	Pentadecane		212	C15H32	000629-62-9	70
4	Nonane		128	C9H20	000111-84-2	60
5	Nonane, 4,5-dimethyl-		156	C11H24	017302-23-7	58



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Sample : M2596-05DL 5X
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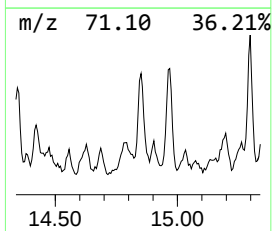
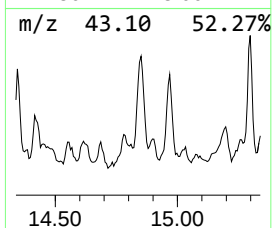
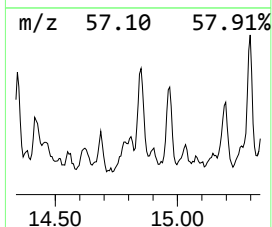
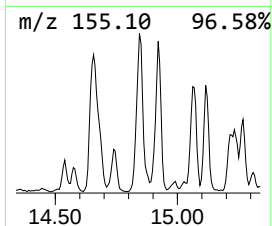
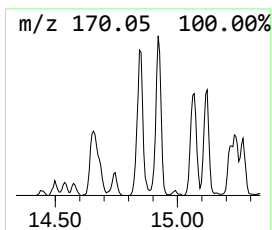
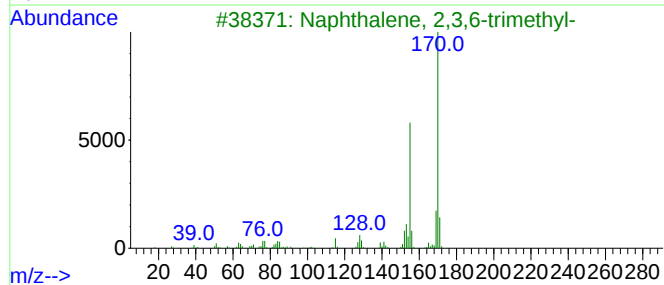
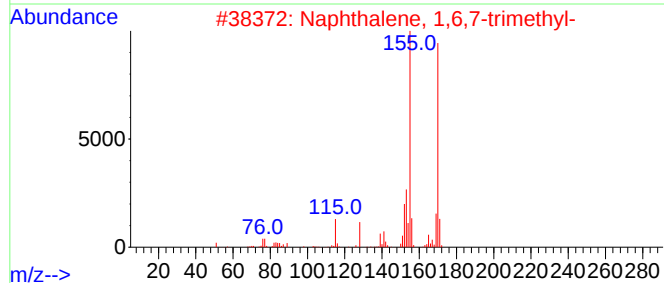
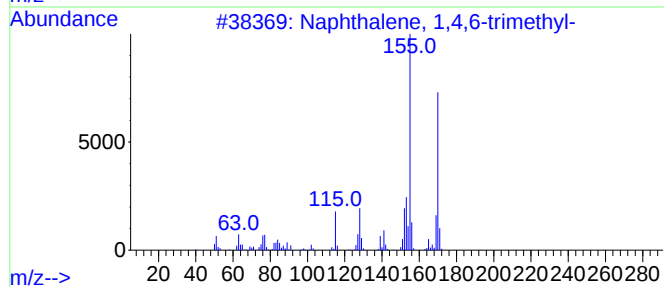
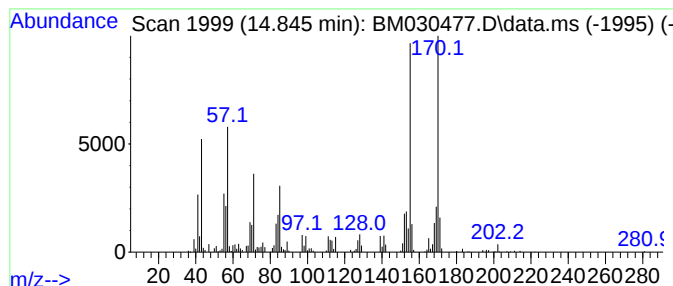
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 1,4,6-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.840	4.50 ng/ul	447545	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	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\
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Sample : M2596-05DL 5X
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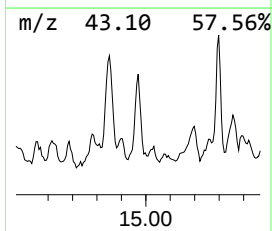
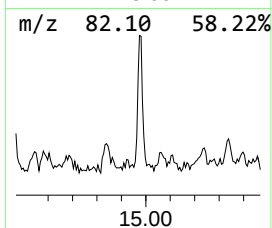
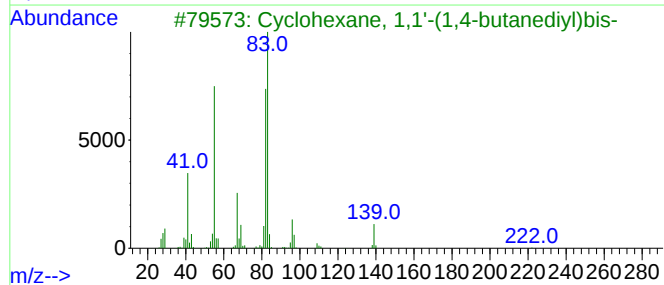
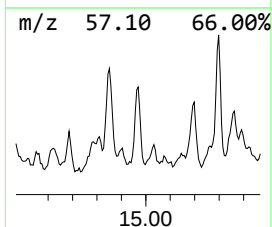
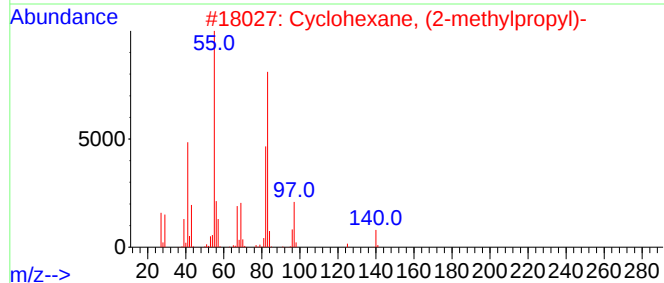
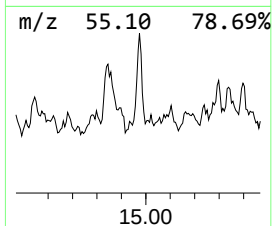
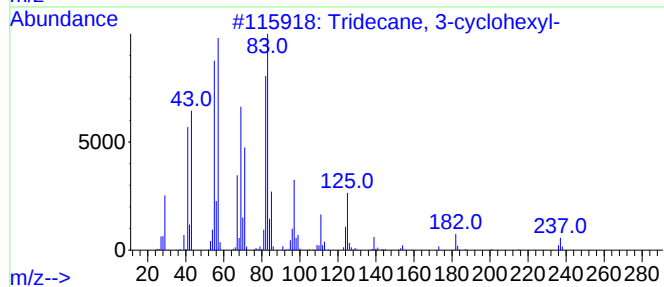
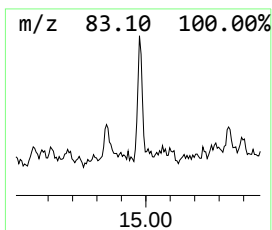
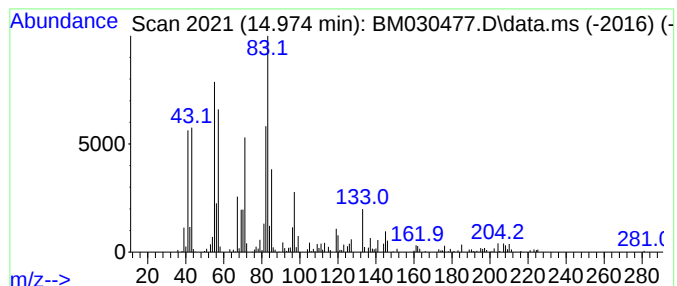
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.970	3.16 ng/ul	314117	Acenaphthene-d10	14.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 3-cyclohexyl-	266	C19H38	013151-88-7	50
2	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	49
3	Cyclohexane, 1,1'-(1,4-butanediyl)-	222	C16H30	006165-44-2	38
4	Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	35
5	Cyclohexane, butyl-	140	C10H20	001678-93-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
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ALS Vial : 15 Sample Multiplier: 1

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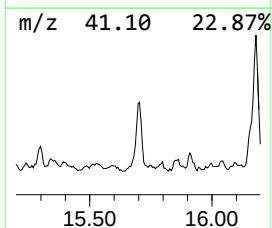
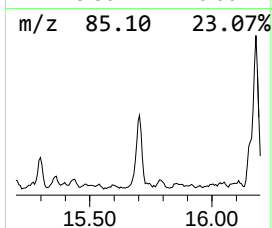
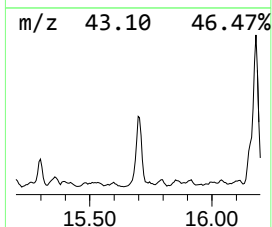
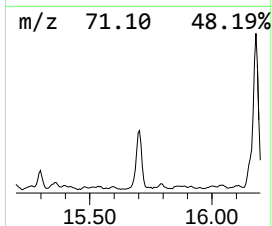
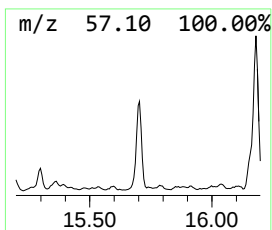
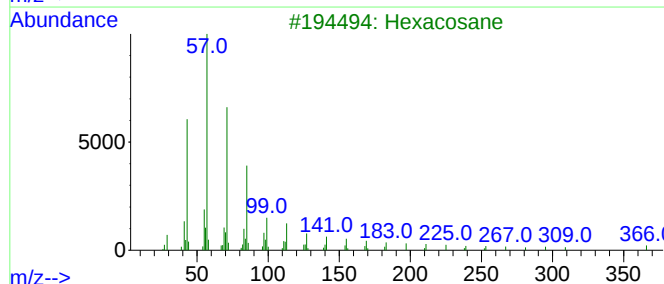
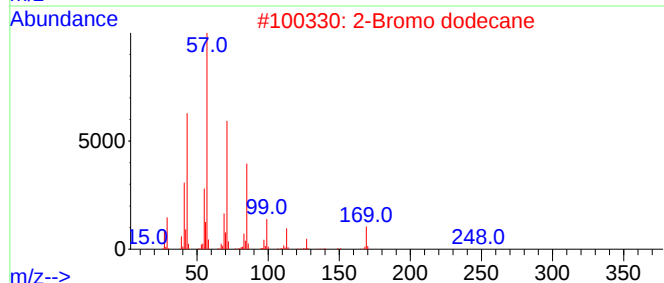
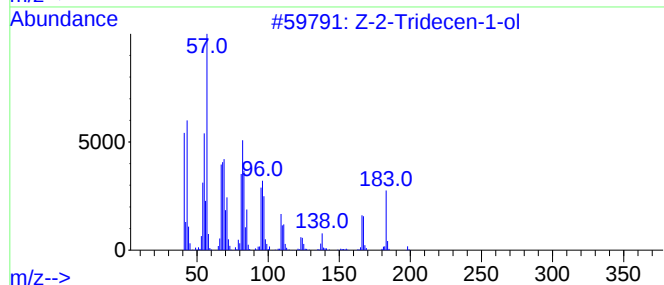
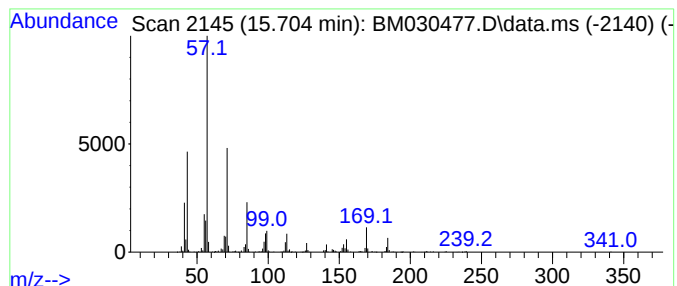
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Z-2-Tridecen-1-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.700	7.71 ng/ul	766841	Acenaphthene-d10	14.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-2-Tridecen-1-ol	198	C13H26O	1000130-75-8	91
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	64
3			Hexacosane	366	C26H54	000630-01-3	58
4			Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	52
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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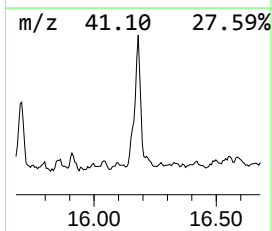
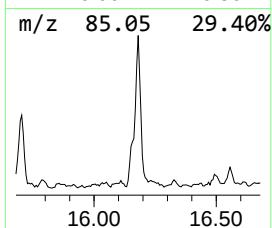
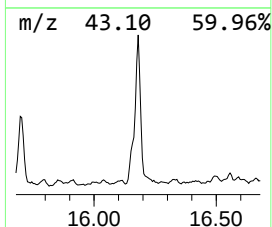
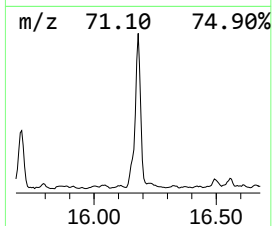
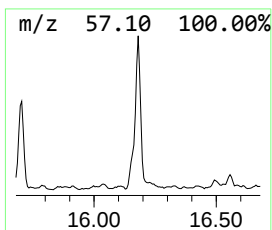
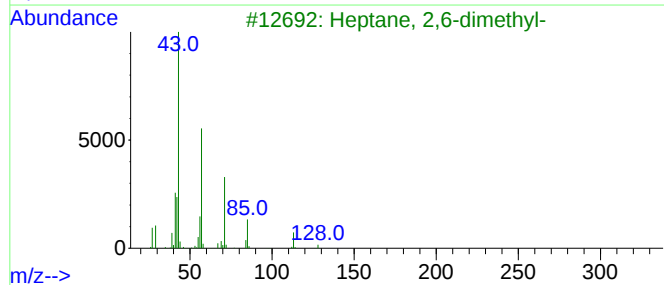
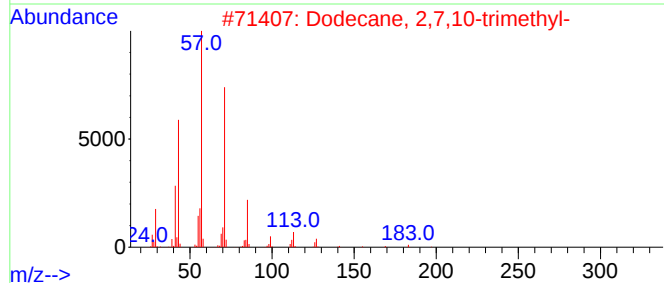
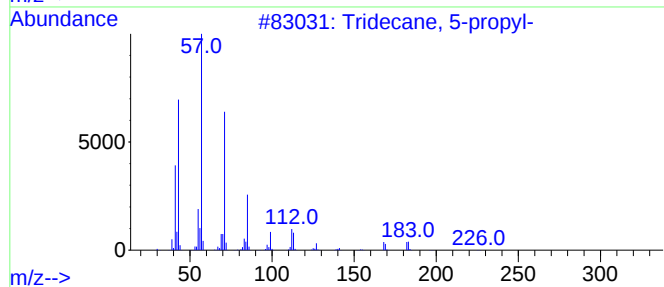
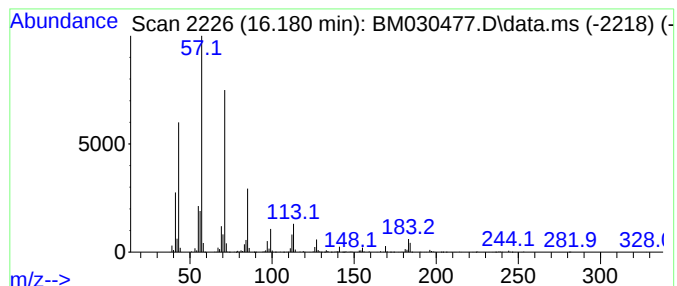
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.180	15.68 ng/ul	1671770	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 5-propyl-	226	C16H34	055045-11-9	93
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
3			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	74
4			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	72
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	72



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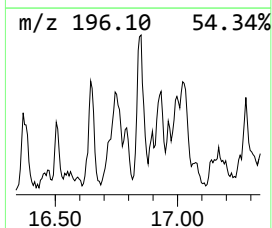
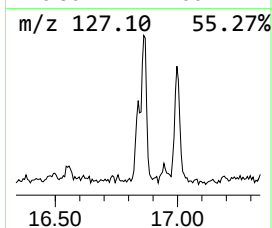
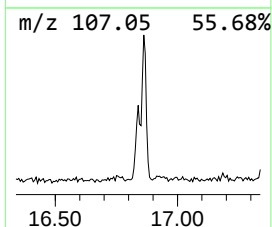
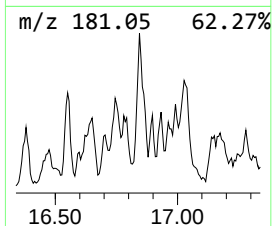
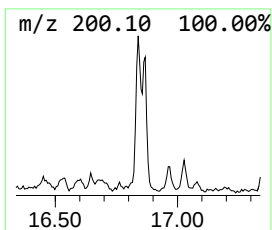
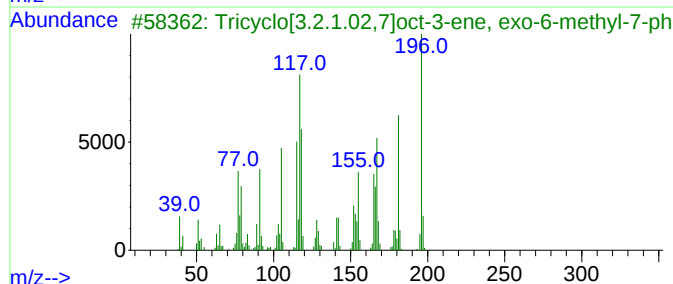
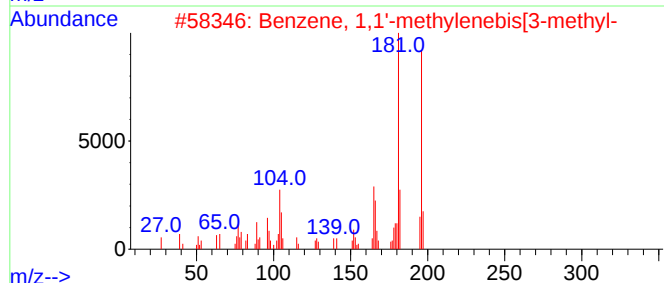
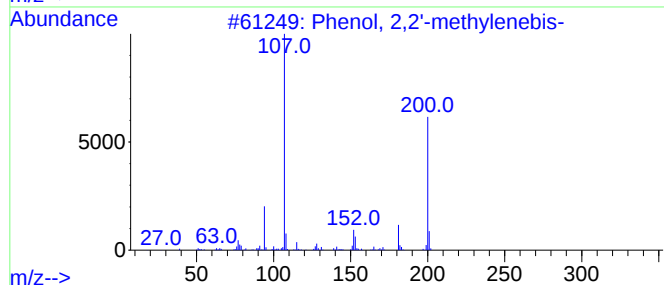
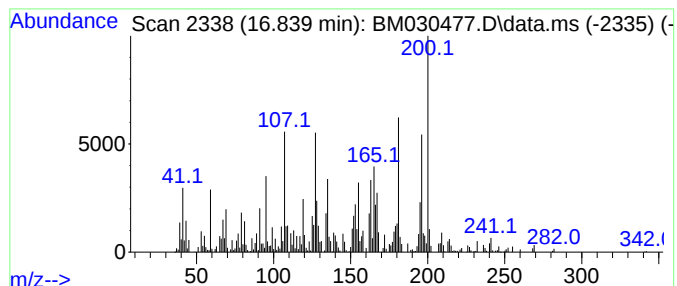
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-01 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.840	2.18 ng/ul	231880	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	42
2			Benzene, 1,1'-methylenebis[3-met...	196	C15H16	021895-14-7	38
3			Tricyclo[3.2.1.0 ^{2,7}]oct-3-ene, e...	196	C15H16	1000142-97-8	35
4			Naphtho[2,1-b]furan, 1,2-dimethyl-	196	C14H12O	129812-23-3	30
5			Benzene, 2,6-dimethyl-1-(phenylm...	196	C15H16	028122-29-4	25



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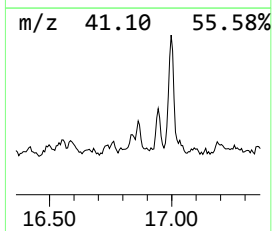
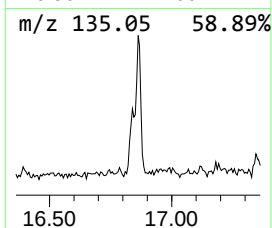
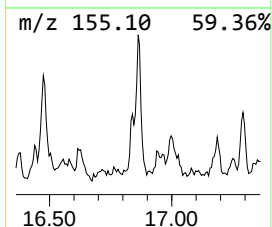
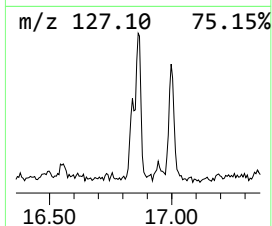
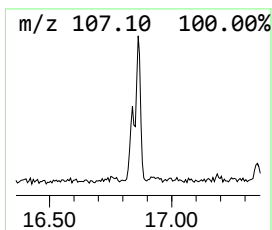
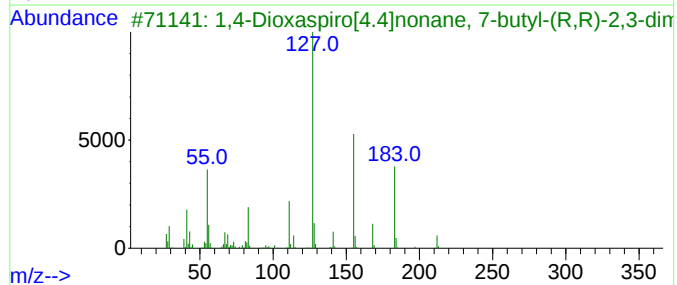
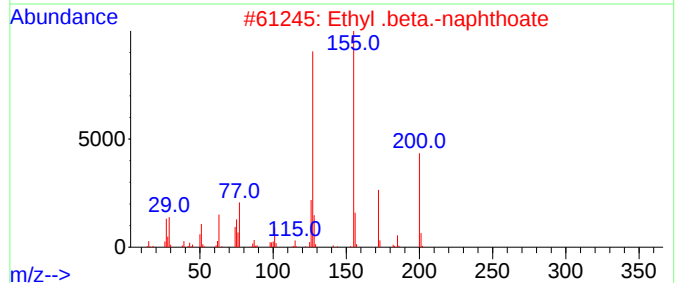
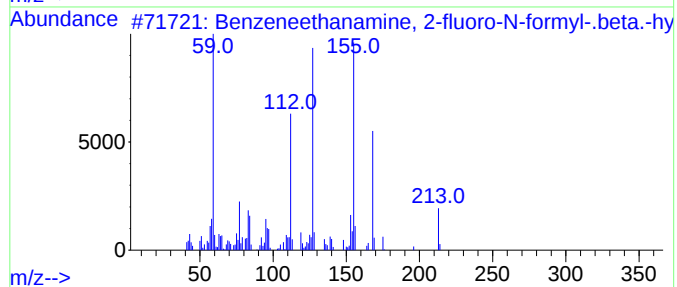
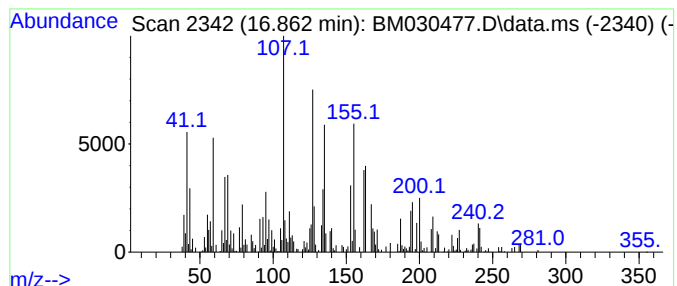
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.860	2.84 ng/ul	302688	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneethanamine, 2-fluoro-N-fo...	213	C10H12FNO3	1000125-03-3	38
2	Ethyl .beta.-naphthoate	200	C13H12O2	003007-91-8	35
3	1,4-Dioxaspiro[4.4]nonane, 7-but...	212	C13H24O2	104807-77-4	25
4	1-Naphthalenecarboxamide, N-(1-m...	227	C15H17NO	031609-18-4	20
5	4-Ethoxy-tricyclo[5.2.1.0(2,6)]d...	201	C13H15NO	1000193-61-5	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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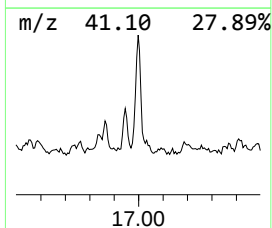
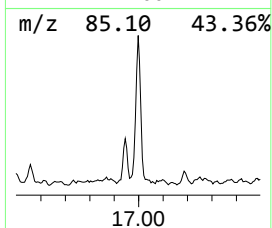
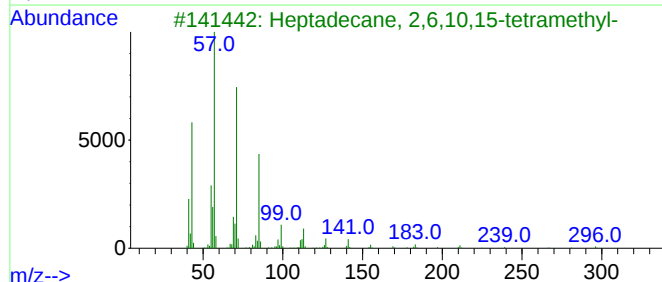
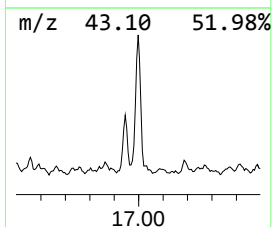
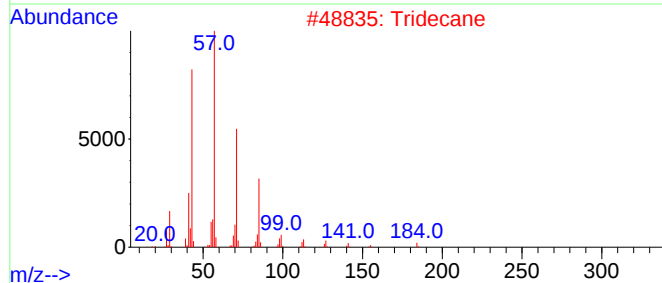
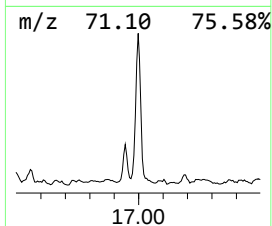
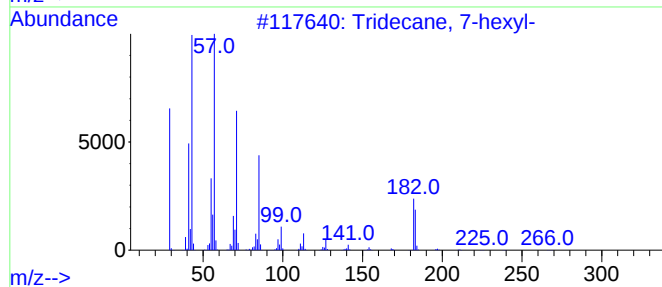
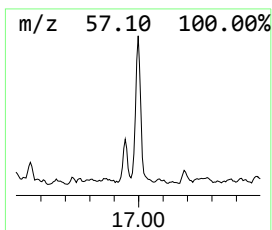
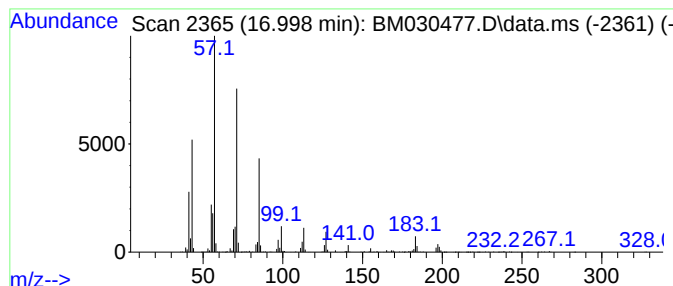
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.000	10.17 ng/ul	1084390	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	94
2		Tridecane	184	C13H28	000629-50-5	91
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
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Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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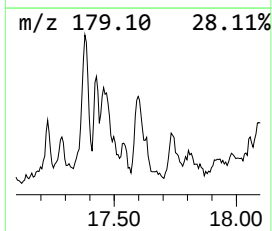
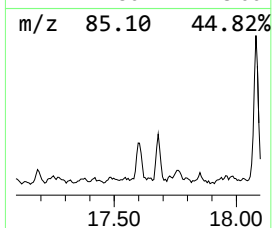
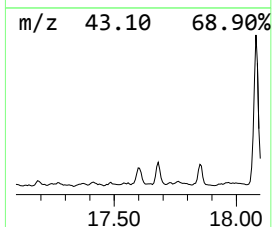
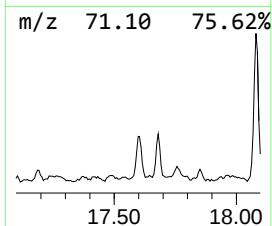
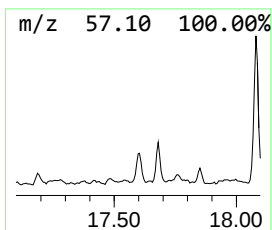
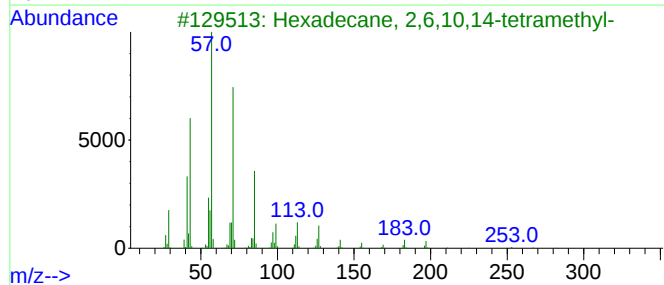
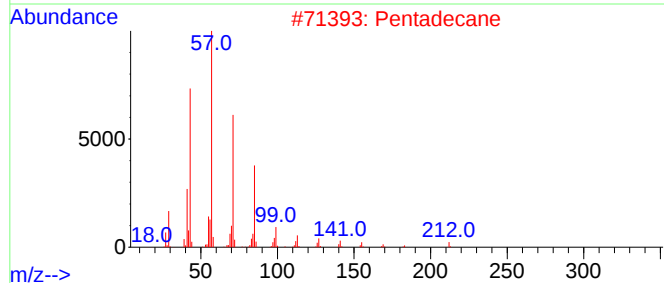
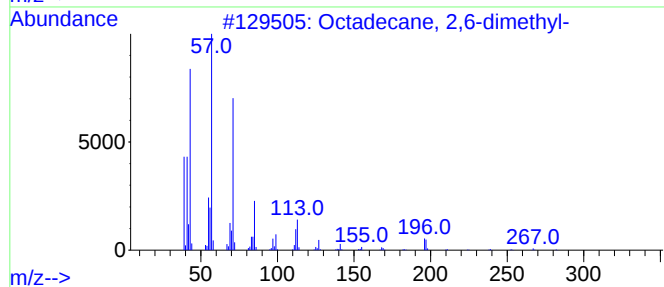
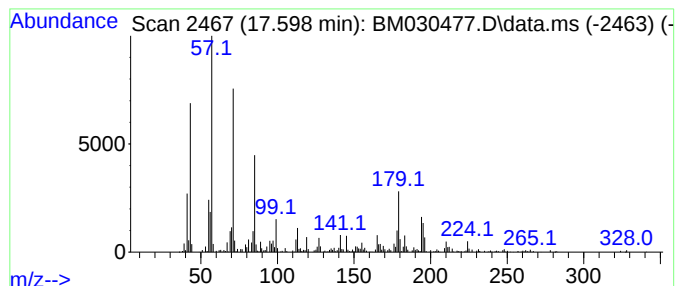
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.600	3.44 ng/ul	366300	Phenanthrene-d10	17.186

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	86
2		Pentadecane	212	C15H32	000629-62-9	83
3		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	64
4		Tridecane	184	C13H28	000629-50-5	64
5		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

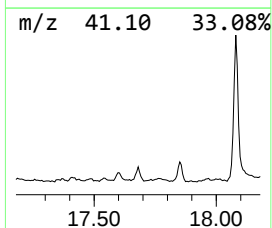
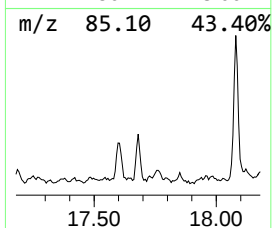
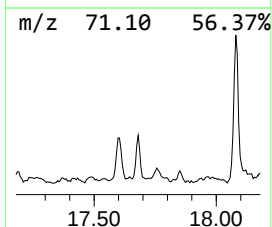
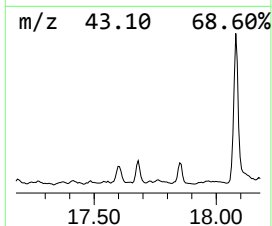
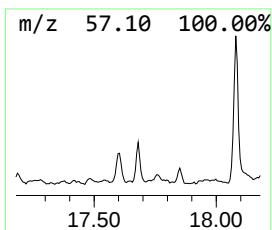
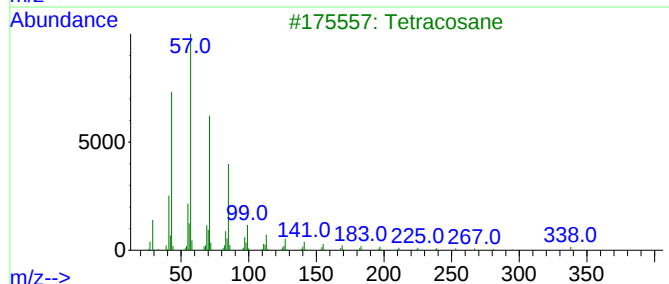
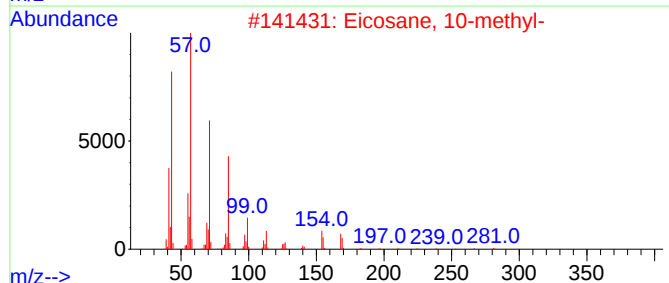
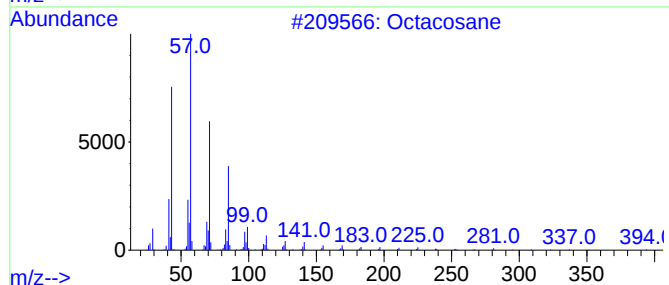
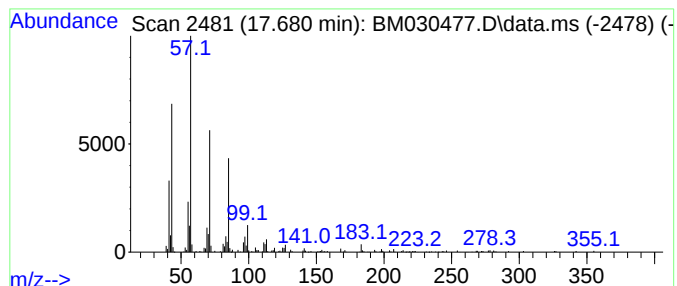
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.680	2.18 ng/ul	232556	Phenanthrene-d10		17.186	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	87
2	Eicosane, 10-methyl-		296	C21H44	054833-23-7	87
3	Tetracosane		338	C24H50	000646-31-1	86
4	Tetradecane		198	C14H30	000629-59-4	86
5	Tetradecane		198	C14H30	000629-59-4	81



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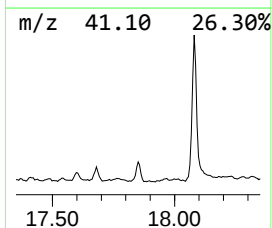
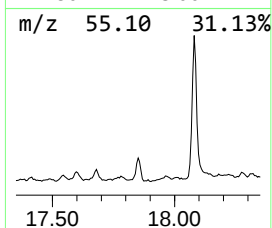
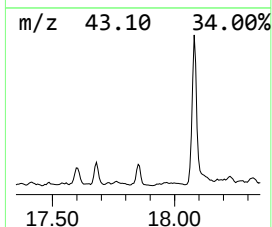
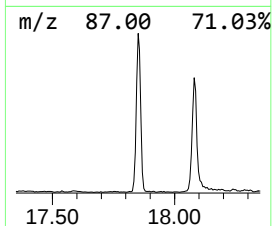
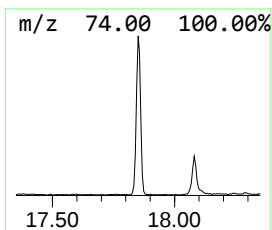
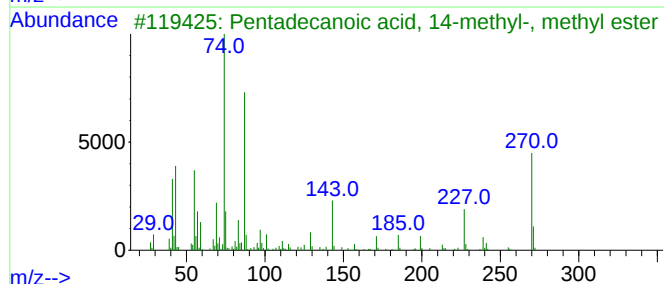
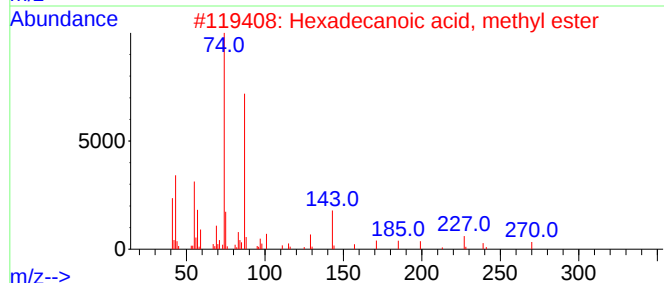
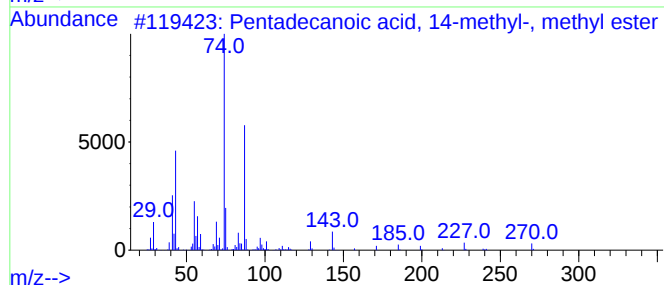
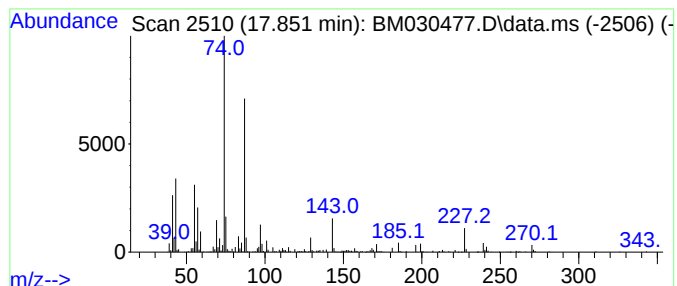
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Pentadecanoic acid, 14-meth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.850	4.47 ng/ul	476274	Phenanthrene-d10	17.186	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
3	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
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Sample : M2596-05DL 5X
Misc :
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ClientSampleId :
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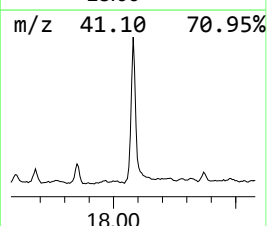
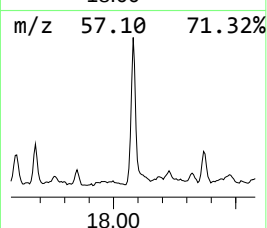
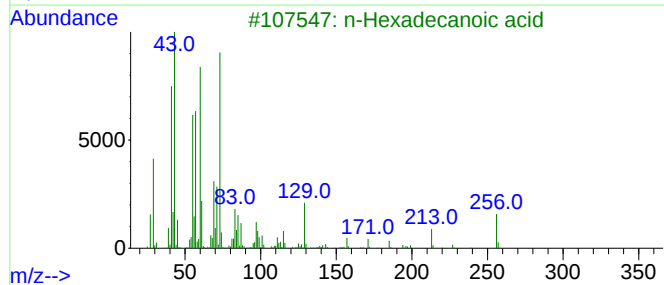
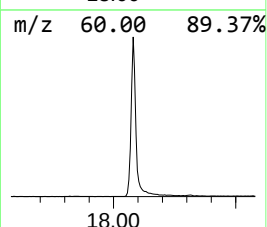
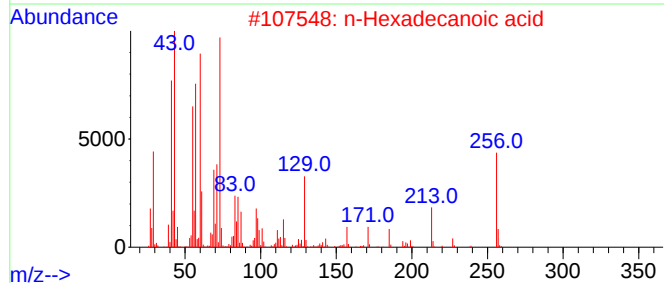
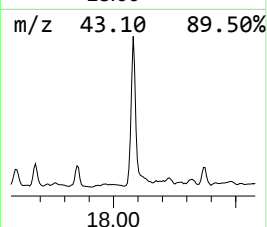
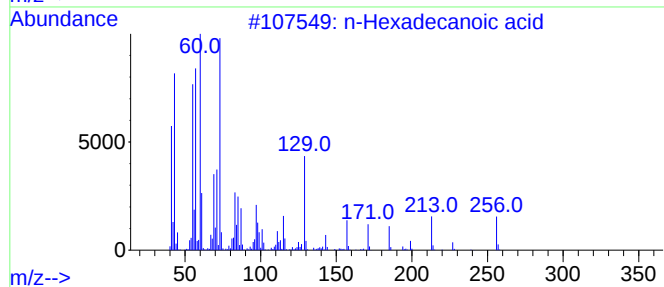
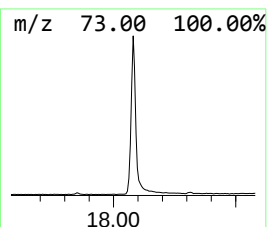
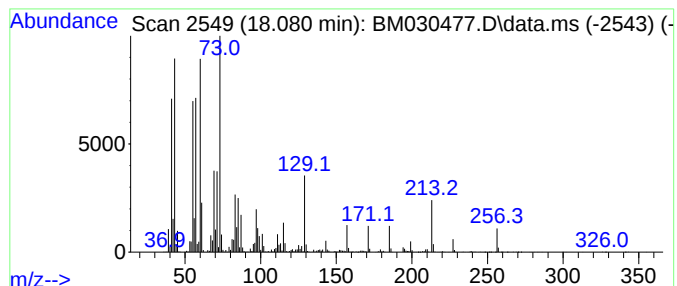
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.080	29.35 ng/ul	3127820	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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	87



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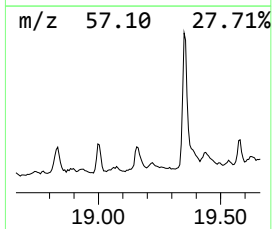
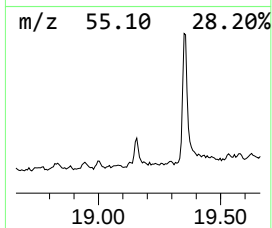
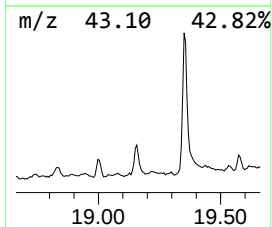
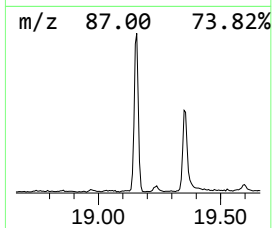
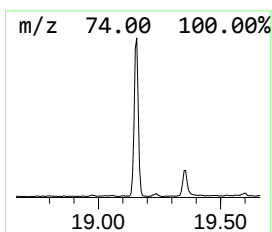
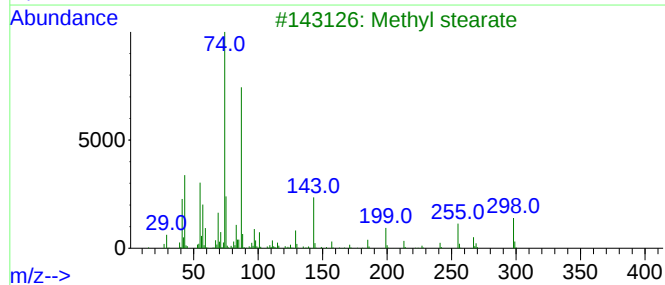
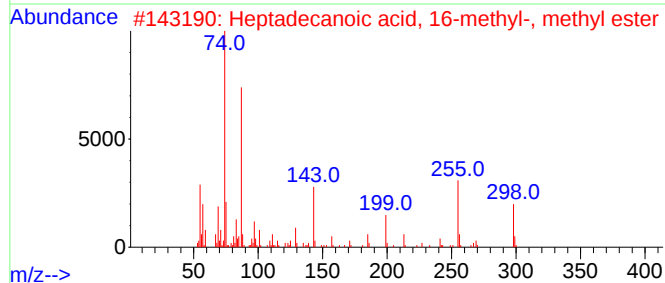
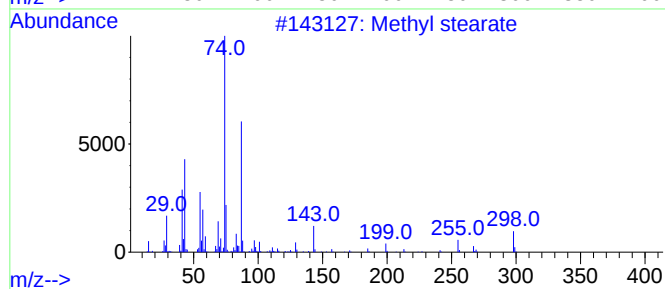
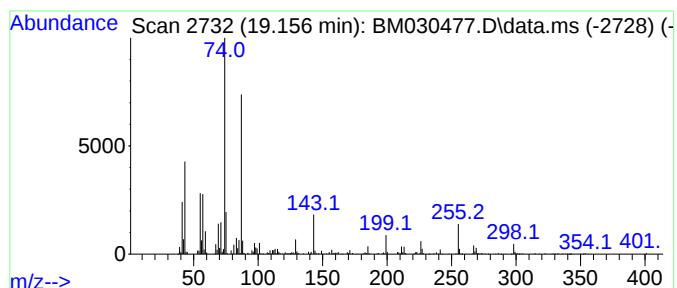
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Methyl stearate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.160	5.09 ng/ul	542364	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	95
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	93
3			Methyl stearate	298	C19H38O2	000112-61-8	87
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	87
5			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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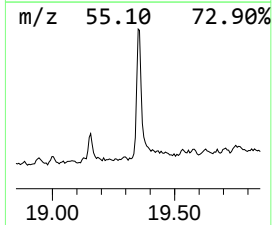
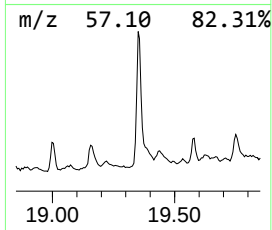
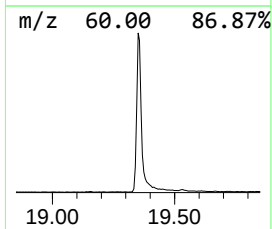
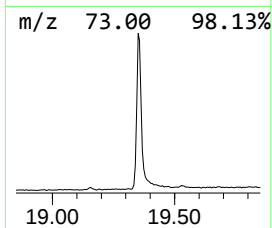
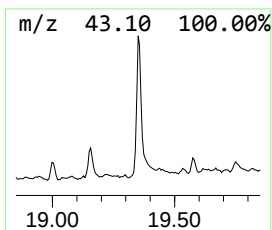
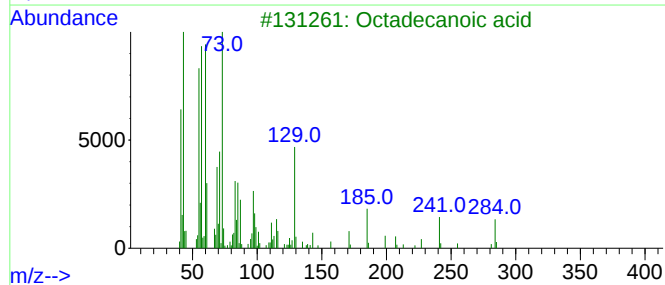
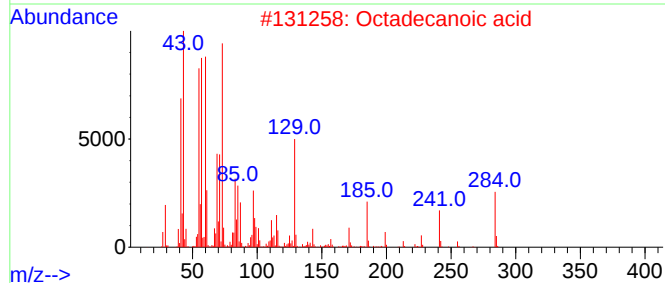
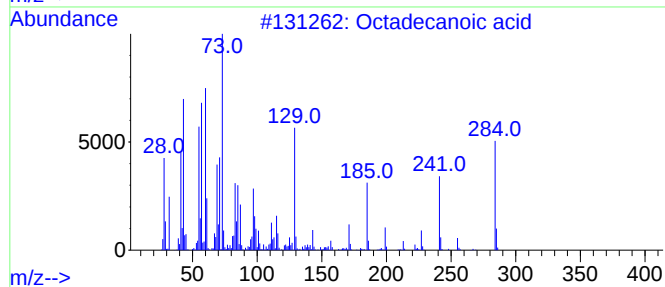
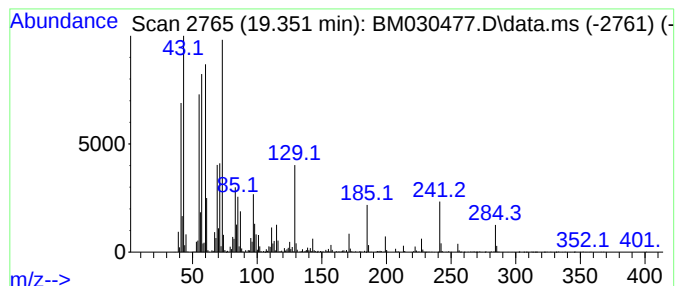
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.350	17.75 ng/ul	2389380	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	97
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5Q1DL

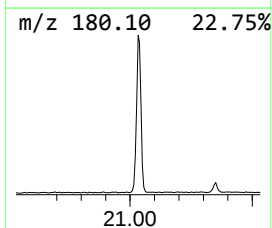
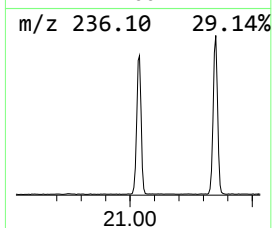
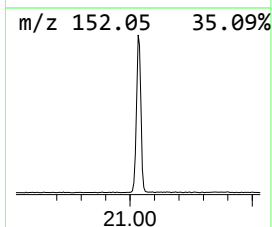
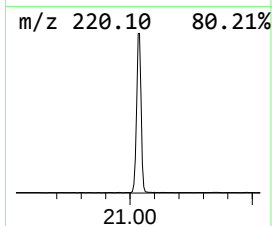
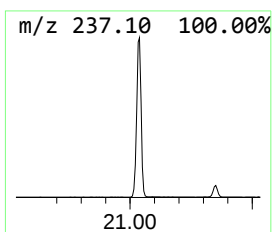
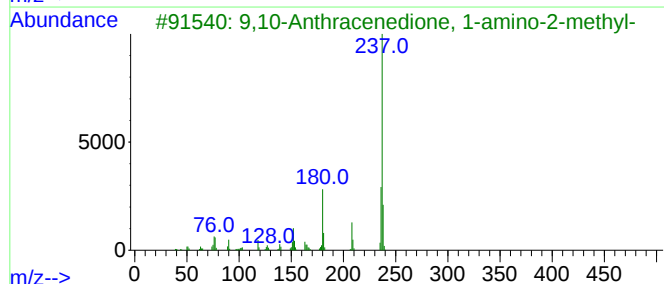
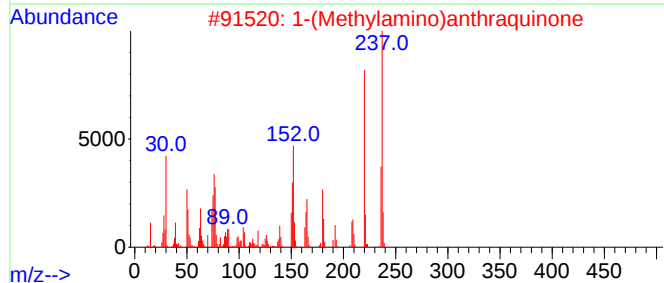
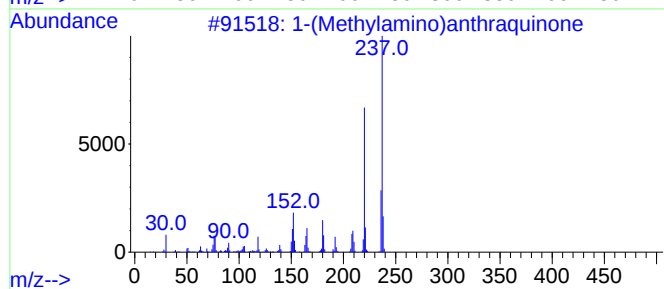
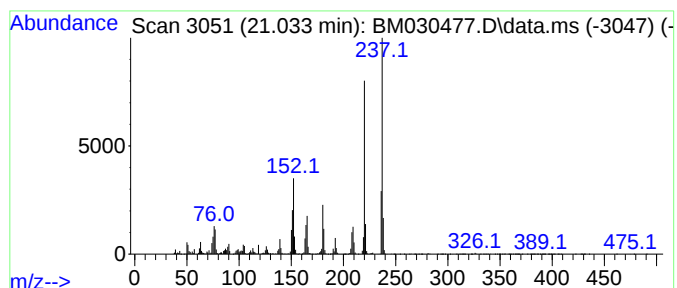
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1-(Methylamino)anthraquinone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.030	35.01 ng/ul	4712460	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	89		
4	2-[p-Methylphenyl]isatogen	237	C15H11NO2	074625-85-7	58		
5	9,10-Anthracenedione, 1-amino-2-...	237	C15H11NO2	000082-28-0	46		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
Data File : BM030477.D
Acq On : 16 Jun 2021 21:46
Operator : CG/JU
Sample : M2596-05DL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG5Q1DL

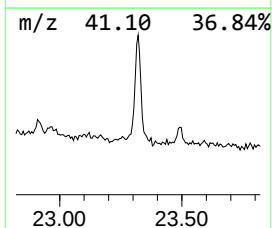
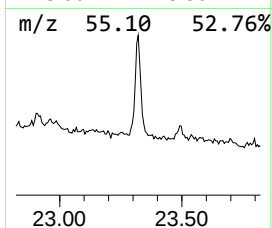
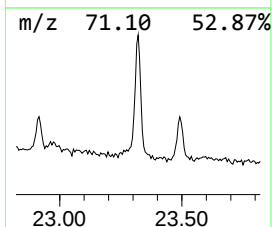
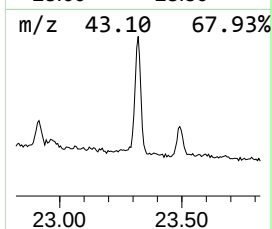
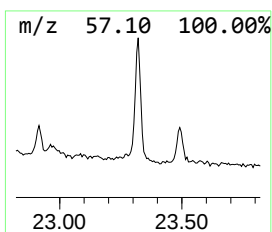
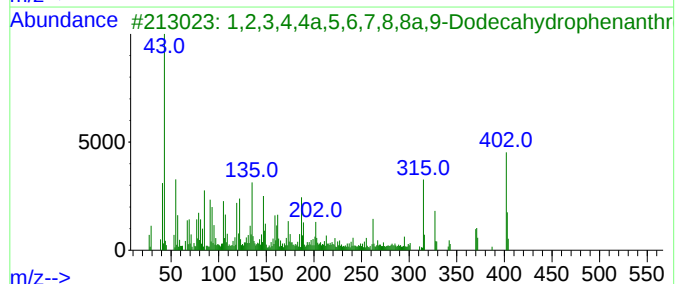
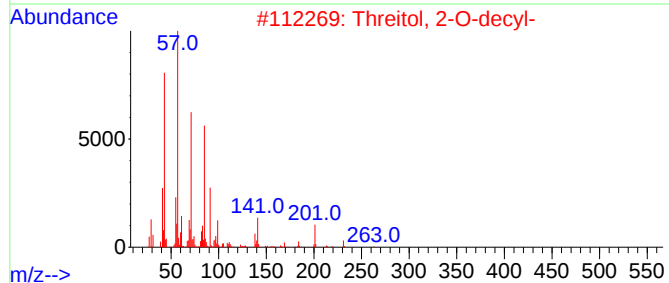
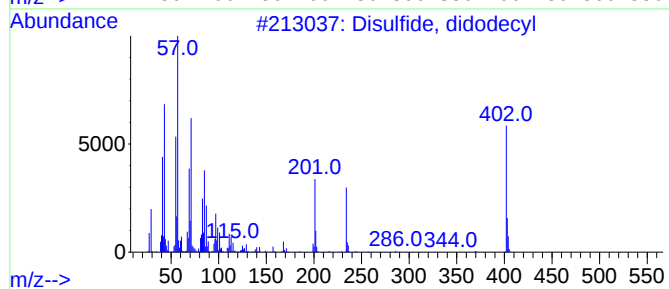
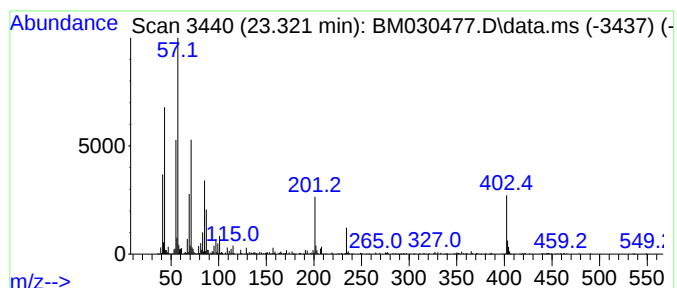
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 Disulfide, didodecyl Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.320	6.48 ng/ul	903139	Perylene-d12	23.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Disulfide, didodecyl	402	C24H50S2	002757-37-1	55
2			Threitol, 2-O-decyl-	262	C14H30O4	1000156-80-6	25
3			1,2,3,4,4a,5,6,7,8,8a,9-Dodecahy...	402	C24H34O5	1000196-44-1	9
4			Dodecane, 1,1'-thiobis-	370	C24H50S	002469-45-6	9
5			Benzenesulfonamide, 3,4-dimethox...	354	C16H22N2O5S	185299-49-4	8



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061621\
 Data File : BM030477.D
 Acq On : 16 Jun 2021 21:46
 Operator : CG/JU
 Sample : M2596-05DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG5Q1DL

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-Butenoic acid...	3.320	30.5	ng/ul	1498050	1	7.828	983251	20.0
2-Propenoic aci...	3.910	2.5	ng/ul	125352	1	7.828	983251	20.0
(DEL) Alkane: S...	10.460	2.9	ng/ul	219160	2	10.610	1489280	20.0
(DEL) Alkane: S...	11.630	4.3	ng/ul	319943	2	10.610	1489280	20.0
(DEL) Alkane: S...	12.250	2.8	ng/ul	209740	2	10.610	1489280	20.0
(DEL) Alkane: S...	12.990	4.3	ng/ul	422162	3	14.451	1988050	20.0
1-Cyclohexene-1...	13.060	2.7	ng/ul	270790	3	14.451	1988050	20.0
1,13-Tetradecad...	13.410	4.0	ng/ul	394426	3	14.451	1988050	20.0
Naphthalene, 1,...	13.610	3.0	ng/ul	297407	3	14.451	1988050	20.0
(DEL) Alkane: C...	14.290	2.1	ng/ul	205089	3	14.451	1988050	20.0
(DEL) Alkane: S...	14.350	2.2	ng/ul	220225	3	14.451	1988050	20.0
Naphthalene, 1,...	14.840	4.5	ng/ul	447545	3	14.451	1988050	20.0
(DEL) Alkane: S...	14.970	3.2	ng/ul	314117	3	14.451	1988050	20.0
Z-2-Tridecen-1-ol	15.700	7.7	ng/ul	766841	3	14.451	1988050	20.0
(DEL) Alkane: S...	16.180	15.7	ng/ul	1671770	4	17.186	2131680	20.0
unknown-01	16.840	2.2	ng/ul	231880	4	17.186	2131680	20.0
unknown-02	16.860	2.8	ng/ul	302688	4	17.186	2131680	20.0
(DEL) Alkane: S...	17.000	10.2	ng/ul	1084390	4	17.186	2131680	20.0
(DEL) Alkane: S...	17.600	3.4	ng/ul	366300	4	17.186	2131680	20.0
(DEL) Alkane: S...	17.680	2.2	ng/ul	232556	4	17.186	2131680	20.0
Pentadecanoic a...	17.850	4.5	ng/ul	476274	4	17.186	2131680	20.0
n-Hexadecanoic ...	18.080	29.4	ng/ul	3127820	4	17.186	2131680	20.0
Methyl stearate	19.160	5.1	ng/ul	542364	4	17.186	2131680	20.0
Octadecanoic acid	19.350	17.8	ng/ul	2389380	5	21.350	2691750	20.0
1-(Methylamino)...	21.030	35.0	ng/ul	4712460	5	21.350	2691750	20.0
Disulfide, dido...	23.320	6.5	ng/ul	903139	6	23.621	2785950	20.0