

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
 Data File : BM043317.D
 Acq On : 14 Dec 2023 00:37
 Operator : MA/JU
 Sample : 05690-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCLH2

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-BM120723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043317.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	47	rVB	111306	172212	1.76%	0.142%
2	3.822	104	108	123	rBV	236800	448630	4.59%	0.369%
3	3.934	123	127	133	rVB	55326	77907	0.80%	0.064%
4	4.051	142	147	153	rVB	72488	115023	1.18%	0.095%
5	4.146	160	163	170	rVB2	24757	38813	0.40%	0.032%
6	4.722	257	261	265	rVB3	28628	41576	0.43%	0.034%
7	5.075	317	321	334	rVB	86226	156655	1.60%	0.129%
8	5.975	469	474	478	rBV3	18914	30118	0.31%	0.025%
9	7.157	668	675	690	rVB	2163061	3767498	38.54%	3.098%
10	7.334	699	705	716	rBV	1630919	2662976	27.24%	2.190%
11	7.528	730	738	745	rBV	2100302	3494155	35.75%	2.873%
12	7.592	745	749	757	rVB2	121504	202960	2.08%	0.167%
13	7.781	775	781	793	rVB	398476	676323	6.92%	0.556%
14	7.998	811	818	827	rBV	1504217	2389847	24.45%	1.965%
15	8.710	931	939	950	rBV	2044523	3335599	34.12%	2.743%
16	8.833	953	960	972	rVB	915750	1516317	15.51%	1.247%
17	8.998	979	988	997	rVB2	37373	100307	1.03%	0.082%
18	9.169	1009	1017	1027	rBV	2046699	3460139	35.40%	2.845%
19	9.263	1030	1033	1038	rVV5	18700	35446	0.36%	0.029%
20	9.551	1078	1082	1088	rVB2	20675	34455	0.35%	0.028%
21	9.892	1126	1140	1152	rBV	1596929	2818367	28.83%	2.317%
22	10.045	1157	1166	1173	rVV	175297	378322	3.87%	0.311%
23	10.110	1175	1177	1186	rVB6	19235	32266	0.33%	0.027%
24	10.251	1194	1201	1207	rVB	93064	153773	1.57%	0.126%
25	10.422	1223	1230	1240	rVV	2470790	4191645	42.88%	3.447%
26	10.804	1287	1295	1302	rBV	2142722	3647236	37.31%	2.999%
27	10.857	1302	1304	1310	rVB	63681	86913	0.89%	0.071%
28	10.951	1311	1320	1341	rVV	1664214	2994193	30.63%	2.462%
29	11.345	1380	1387	1394	rBV2	89400	163858	1.68%	0.135%
30	11.598	1424	1430	1441	rBV2	44236	100747	1.03%	0.083%
31	11.839	1464	1471	1476	rVB5	18464	37174	0.38%	0.031%
32	12.174	1519	1528	1533	rBV	42629	83289	0.85%	0.068%
33	12.386	1553	1564	1569	rVV3	47574	98012	1.00%	0.081%
34	12.463	1571	1577	1582	rVB	108799	169506	1.73%	0.139%
35	12.680	1604	1614	1620	rVB	90041	170161	1.74%	0.140%
36	13.457	1742	1746	1751	rVV2	49184	88755	0.91%	0.073%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
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37	13.527	1751	1758	1767	rVB2	62243	113638	1.16%	0.093%
38	13.792	1796	1803	1804	rBV3	32310	48649	0.50%	0.040%
39	13.810	1804	1806	1813	rVB7	33942	45651	0.47%	0.038%
40	13.898	1816	1821	1824	rBV6	21361	28556	0.29%	0.023%
41	13.951	1824	1830	1835	rBV3	67249	119725	1.22%	0.098%
42	14.045	1835	1846	1853	rVV	4239325	5894064	60.30%	4.846%
43	14.110	1853	1857	1861	rVB2	52683	72930	0.75%	0.060%
44	14.227	1873	1877	1882	rVB5	50274	70523	0.72%	0.058%
45	14.321	1885	1893	1902	rBV	5030367	7402633	75.73%	6.087%
46	14.463	1911	1917	1921	rBV2	121971	184752	1.89%	0.152%
47	14.510	1921	1925	1935	rVB3	117258	185311	1.90%	0.152%
48	14.627	1935	1945	1950	rBV	2912749	4631668	47.38%	3.808%
49	14.686	1950	1955	1962	rVB	729484	1045434	10.70%	0.860%
50	14.821	1971	1978	1983	rBV	1986374	2859758	29.26%	2.351%
51	14.868	1983	1986	1993	rVV	114508	175459	1.80%	0.144%
52	14.939	1993	1998	2001	rVV2	64498	95845	0.98%	0.079%
53	14.980	2001	2005	2008	rVV	117080	163061	1.67%	0.134%
54	15.021	2008	2012	2021	rVB	308793	603642	6.18%	0.496%
55	15.286	2053	2057	2061	rVV6	30681	51357	0.53%	0.042%
56	15.415	2072	2079	2082	rBV5	16285	36565	0.37%	0.030%
57	15.615	2105	2113	2118	rBV	6116449	8638637	88.38%	7.103%
58	15.668	2118	2122	2128	rVB	433976	624685	6.39%	0.514%
59	15.733	2128	2133	2140	rBV	575166	791454	8.10%	0.651%
60	15.833	2146	2150	2151	rBV2	26533	29909	0.31%	0.025%
61	15.957	2167	2171	2175	rBV	48157	67284	0.69%	0.055%
62	16.080	2186	2192	2203	rVB3	83784	249330	2.55%	0.205%
63	16.186	2203	2210	2216	rBV3	93005	155570	1.59%	0.128%
64	16.333	2229	2235	2242	rBV4	79470	138424	1.42%	0.114%
65	16.668	2282	2292	2296	rBV6	43753	104560	1.07%	0.086%
66	16.827	2313	2319	2323	rBV2	35077	62843	0.64%	0.052%
67	16.909	2328	2333	2341	rVB3	86576	143438	1.47%	0.118%
68	17.021	2348	2352	2360	rVB4	26097	48506	0.50%	0.040%
69	17.121	2367	2369	2372	rVV4	23570	29151	0.30%	0.024%
70	17.180	2375	2379	2387	rVV	78978	127661	1.31%	0.105%
71	17.362	2401	2410	2414	rBV	3325734	4871470	49.84%	4.006%
72	17.404	2414	2417	2422	rVB	866926	1199842	12.28%	0.987%
73	17.462	2422	2427	2432	rBV2	5952949	8563066	87.60%	7.041%
74	17.562	2439	2444	2449	rVB3	92064	129653	1.33%	0.107%
75	17.768	2473	2479	2485	rVB	130388	213684	2.19%	0.176%
76	17.868	2492	2496	2504	rBV2	39820	74793	0.77%	0.061%

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77	18.039	2521	2525	2528	rVB5	25259	32409	0.33%	0.027%
78	18.139	2537	2542	2550	rVB	46070	78101	0.80%	0.064%
79	18.262	2555	2563	2573	rBV2	759302	1232903	12.61%	1.014%
80	18.433	2587	2592	2596	rBV	183008	289589	2.96%	0.238%
81	18.586	2616	2618	2624	rVB4	24226	34070	0.35%	0.028%
82	18.739	2641	2644	2648	rVB	51105	65212	0.67%	0.054%
83	18.839	2657	2661	2668	rVB	165937	224819	2.30%	0.185%
84	18.980	2682	2685	2690	rVB2	53244	60817	0.62%	0.050%
85	19.103	2702	2706	2711	rBV	119097	167068	1.71%	0.137%
86	19.156	2711	2715	2718	rBV6	30909	45477	0.47%	0.037%
87	19.315	2739	2742	2751	rVB4	129329	211088	2.16%	0.174%
88	19.415	2755	2759	2764	rVB	748697	992062	10.15%	0.816%
89	19.474	2764	2769	2773	rVB2	253396	342135	3.50%	0.281%
90	19.533	2776	2779	2788	rBV	201599	291441	2.98%	0.240%
91	19.750	2808	2816	2829	rVB	6787673	9774656	100.00%	8.037%
92	20.268	2901	2904	2907	rBV2	78732	90399	0.92%	0.074%
93	21.250	3067	3071	3081	rBV	1402597	1869084	19.12%	1.537%
94	21.533	3114	3119	3124	rBV	2937946	3819531	39.08%	3.141%
95	21.709	3146	3149	3154	rVB	145709	171189	1.75%	0.141%
96	22.191	3226	3231	3243	rVB	1261761	1791720	18.33%	1.473%
97	23.297	3415	3419	3429	rVB	385918	661369	6.77%	0.544%
98	23.797	3497	3504	3514	rVB2	3070362	5998306	61.37%	4.932%
99	23.956	3524	3531	3538	rVB	1533437	3160753	32.34%	2.599%
100	27.638	4151	4157	4169	rVB4	669520	2149679	21.99%	1.768%

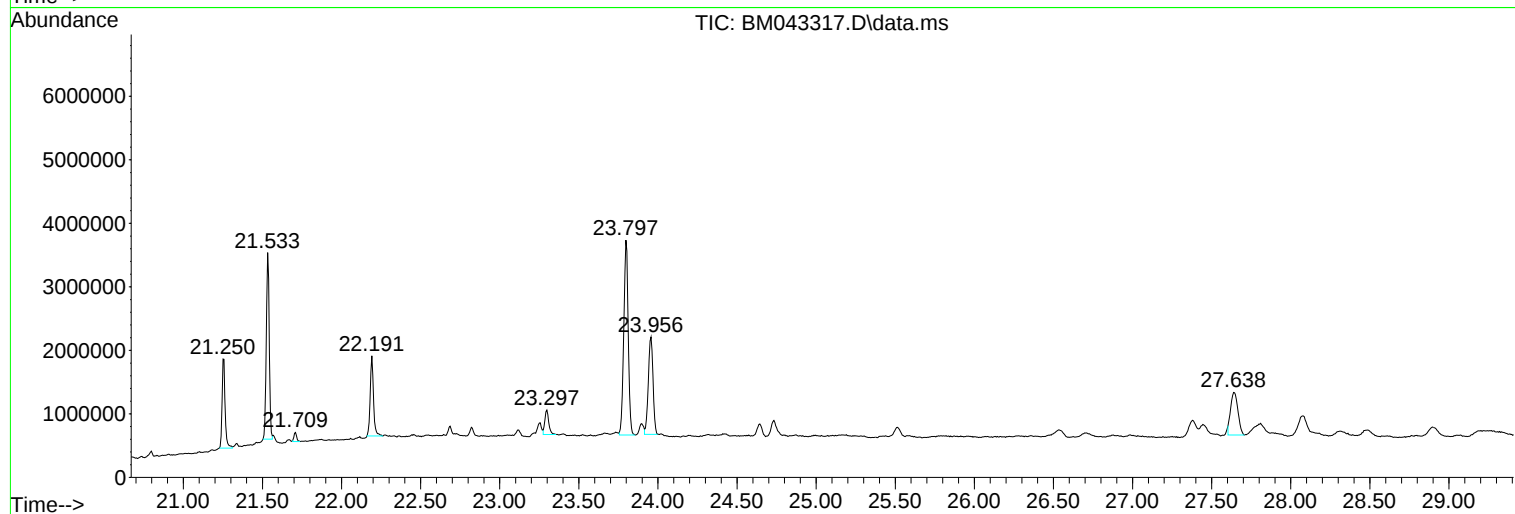
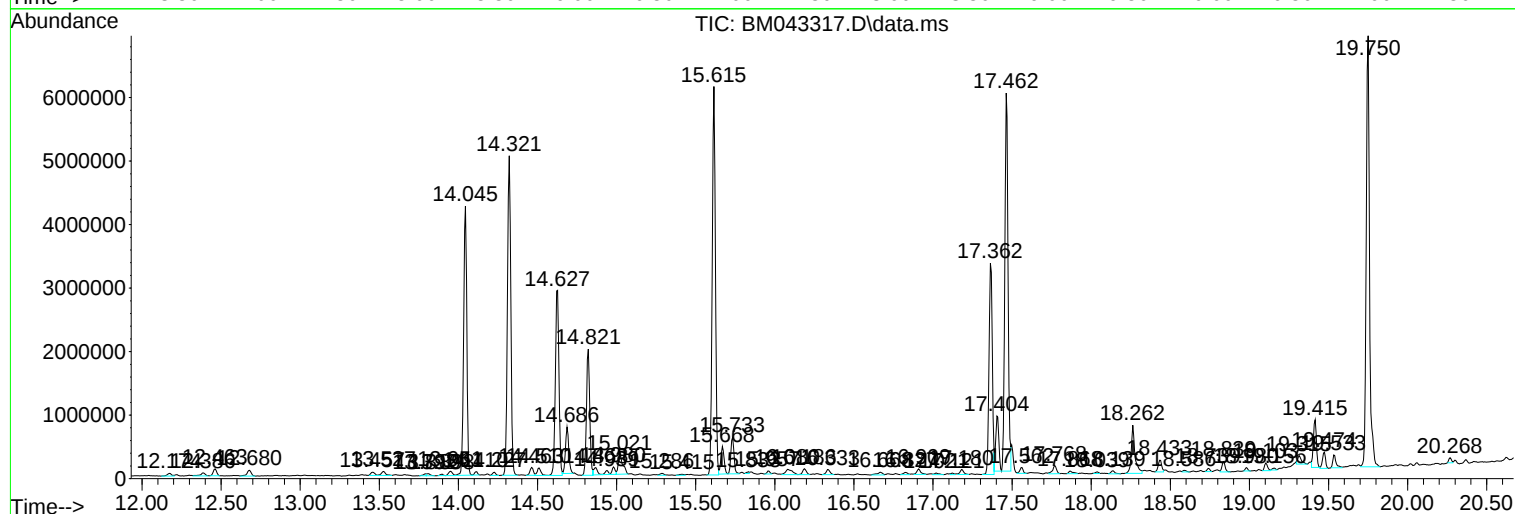
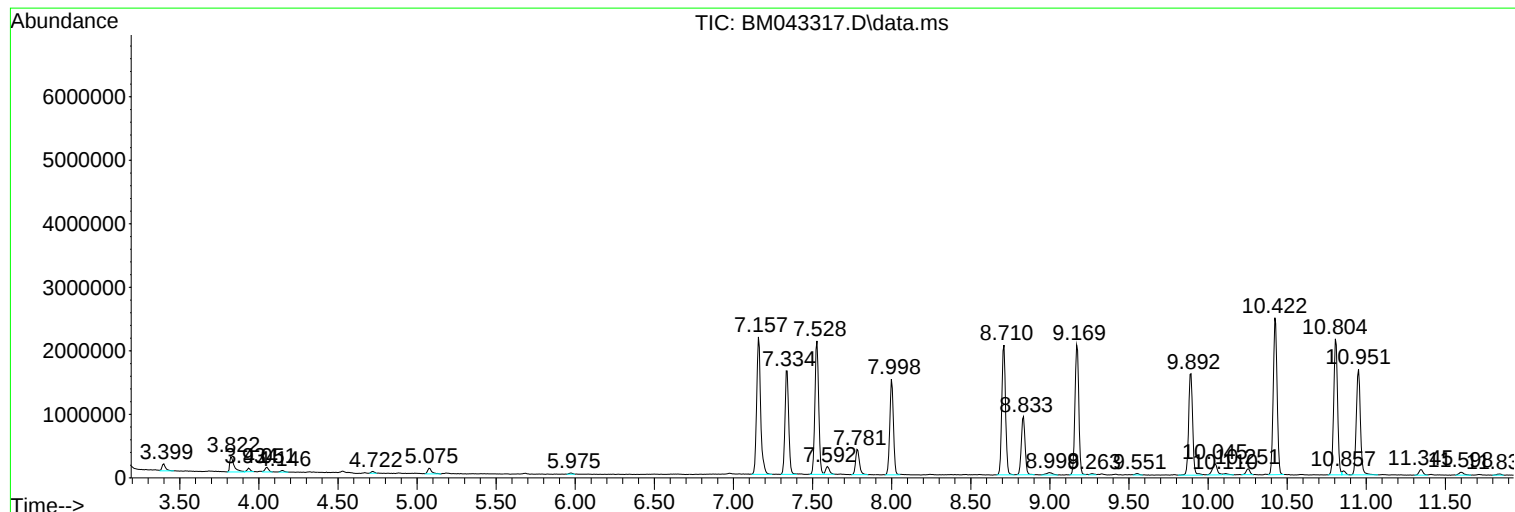
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Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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



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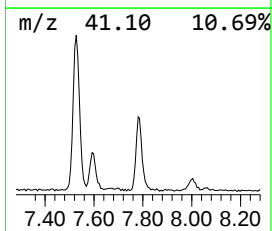
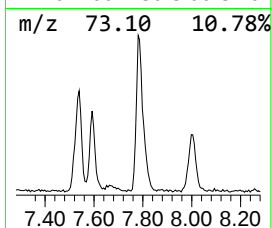
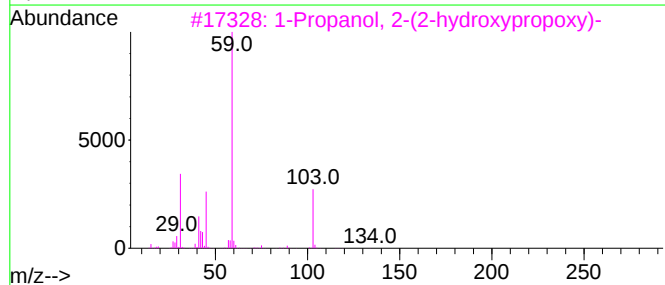
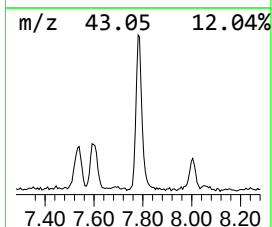
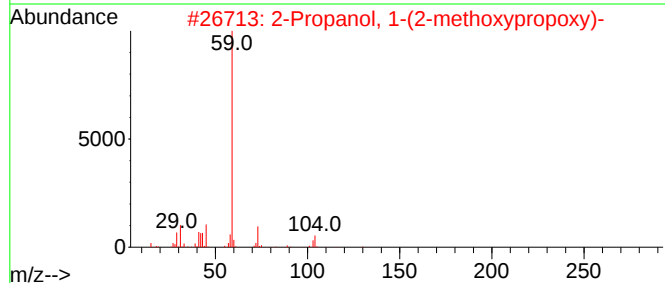
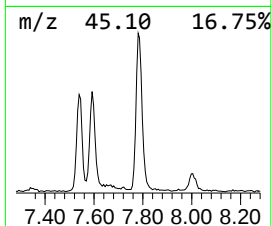
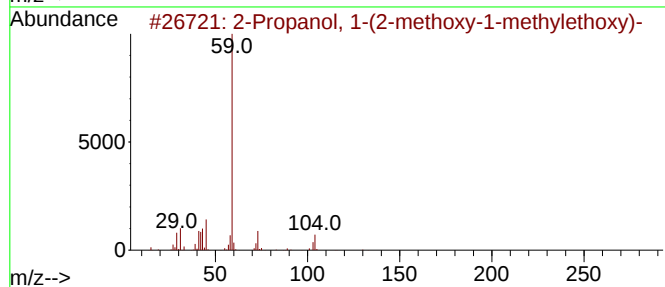
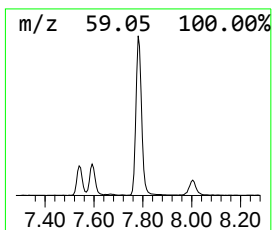
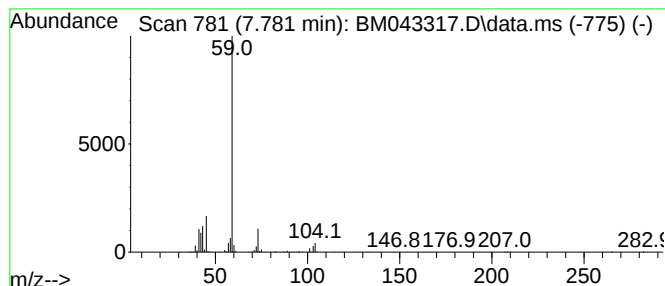
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.781	5.66 ng/ul	676323	1,4-Dichlorobenzene-d4	7.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	74		
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	74		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		
4	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	64		
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64		



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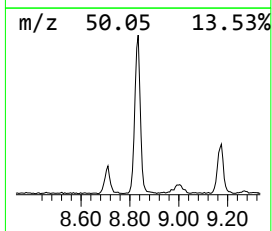
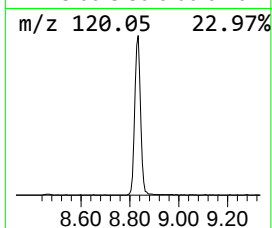
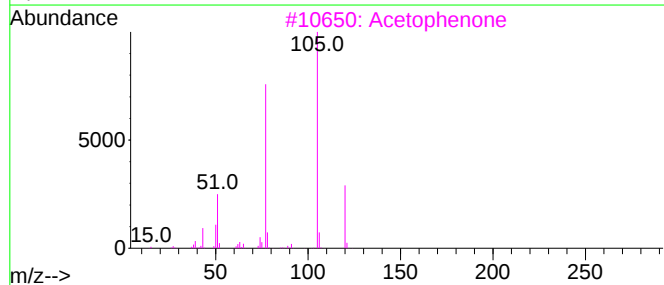
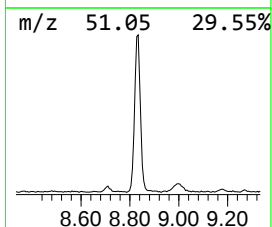
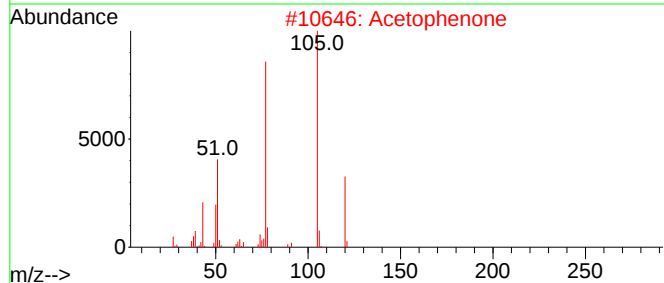
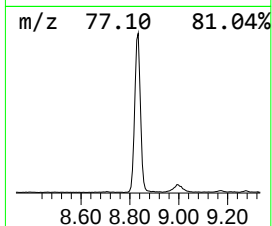
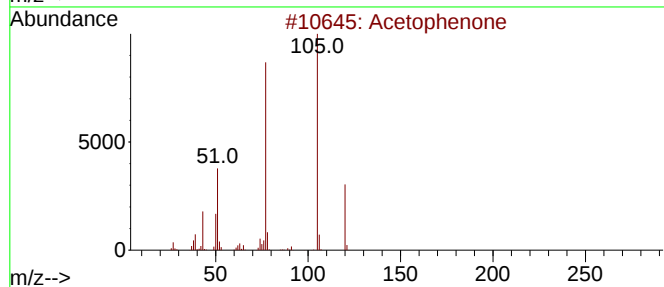
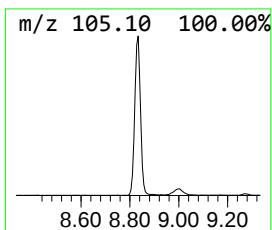
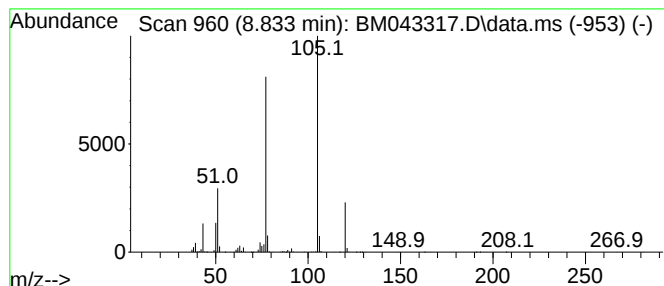
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Aceto phenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.833	12.69 ng/ul	1516320	1,4-Dichlorobenzene-d4	7.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	94
2	Acetophenone			120	C8H8O	000098-86-2	94
3	Acetophenone			120	C8H8O	000098-86-2	93
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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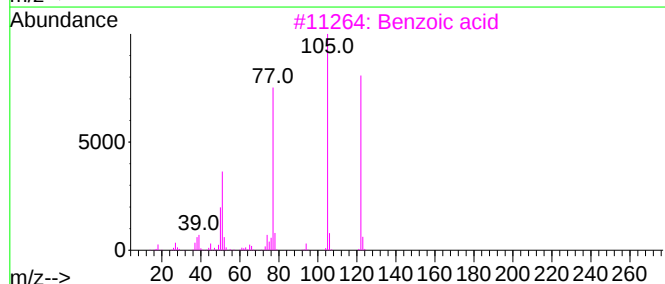
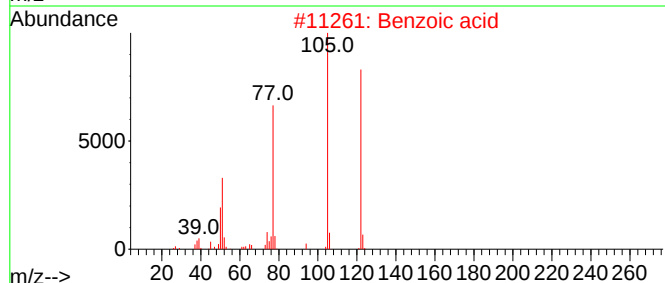
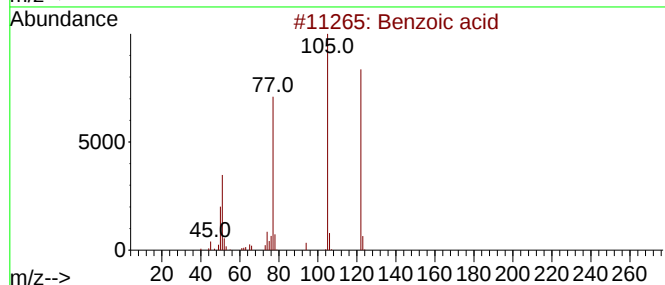
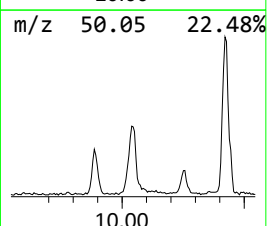
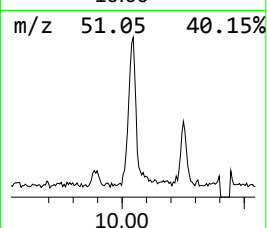
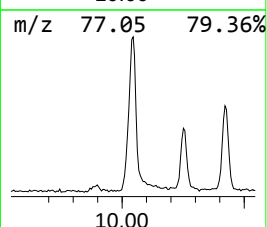
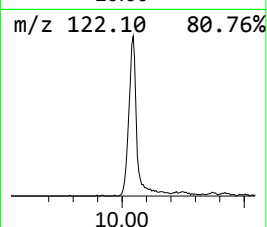
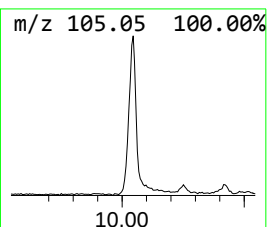
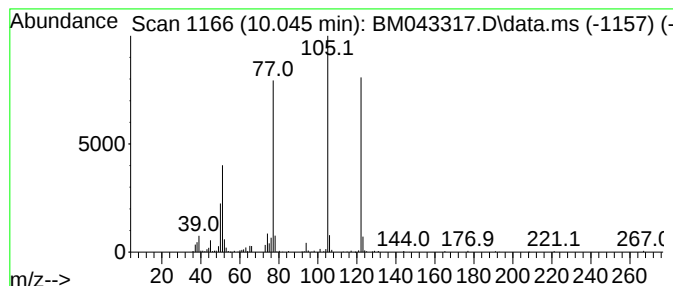
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.045	2.07 ng/ul	378322	Naphthalene-d8	10.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	96
2			Benzoic acid	122	C7H6O2	000065-85-0	94
3			Benzoic acid	122	C7H6O2	000065-85-0	94
4			Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	91
5			Benzoic acid	122	C7H6O2	000065-85-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
Operator : MA/JU
Sample : 05690-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLH2

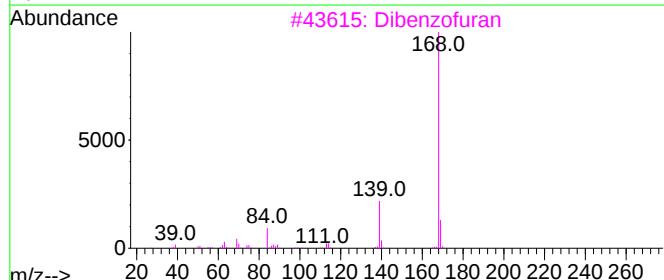
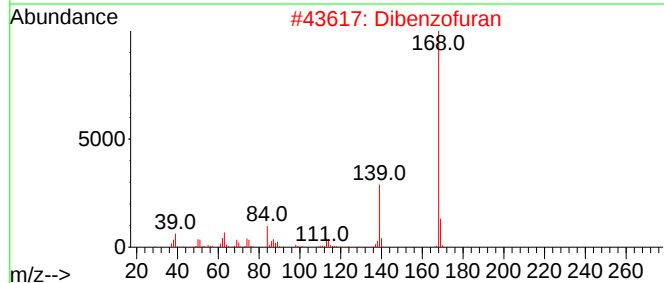
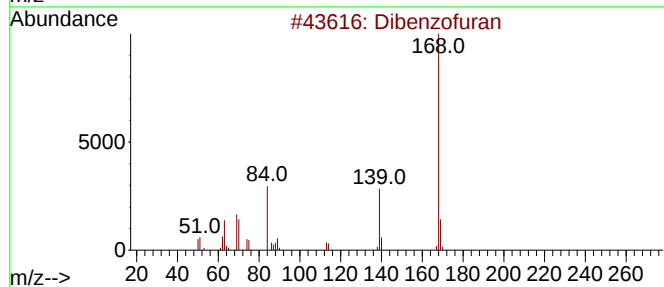
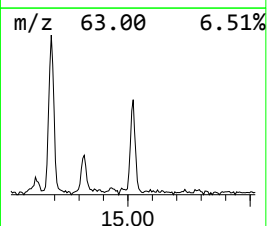
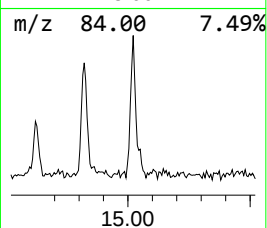
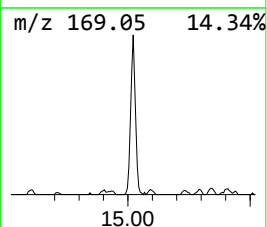
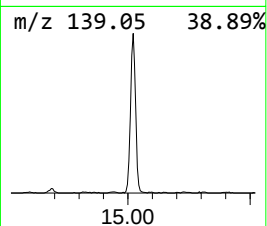
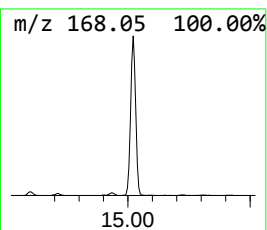
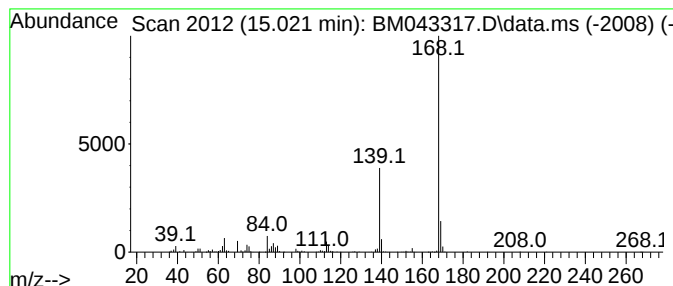
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dibenzo furan Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.021	2.61 ng/ul	603642	Acenaphthene-d10	14.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	86
2			Dibenzofuran	168	C12H8O	000132-64-9	86
3			Dibenzofuran	168	C12H8O	000132-64-9	59
4			3,5-Dimethoxybenzyl alcohol	168	C9H12O3	000705-76-0	56
5			1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
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Sample : 05690-03
Misc :
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Instrument :
BNA_M
ClientSampleId :
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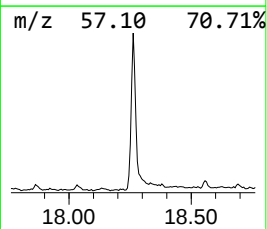
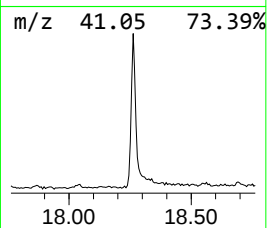
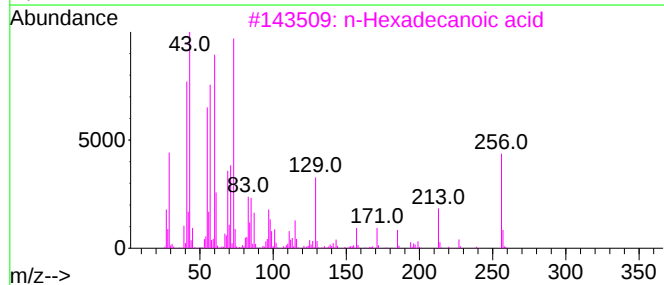
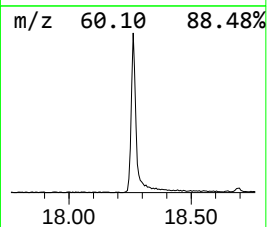
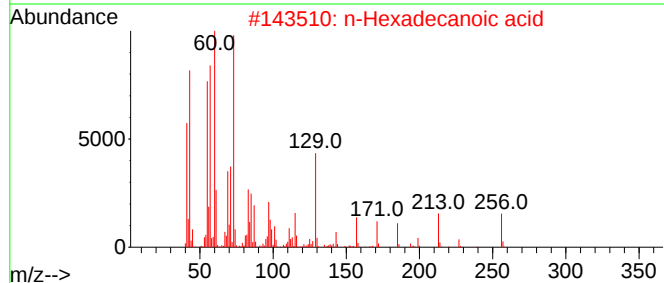
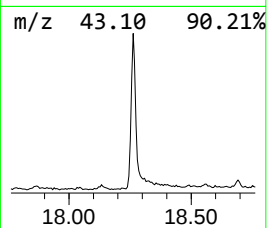
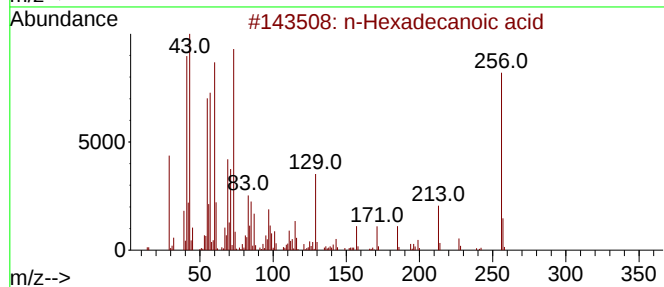
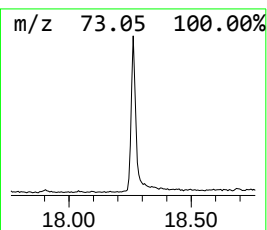
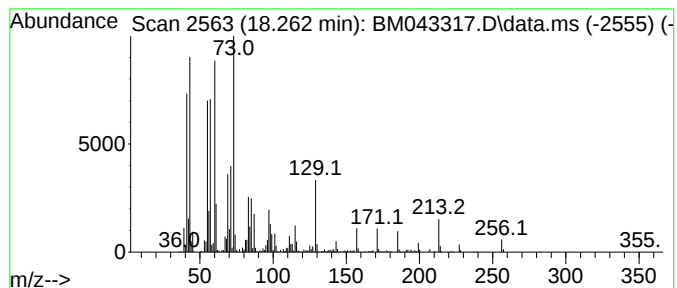
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.262	5.06 ng/ul	1232900	Phenanthrene-d10	17.362

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
Operator : MA/JU
Sample : 05690-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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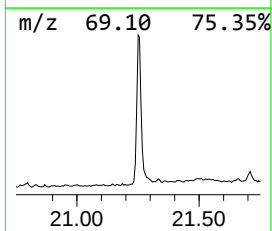
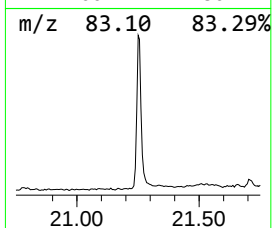
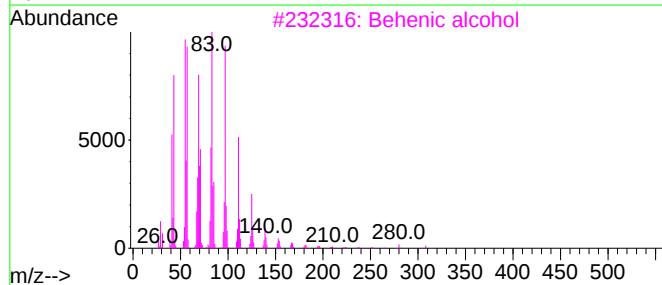
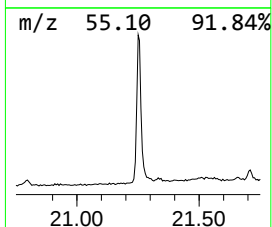
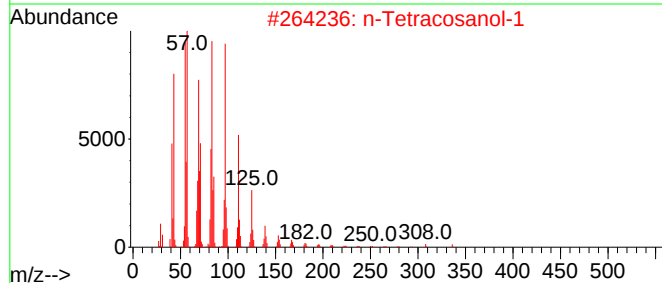
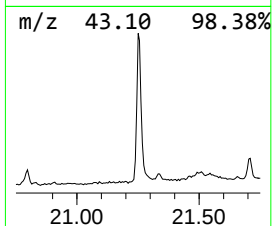
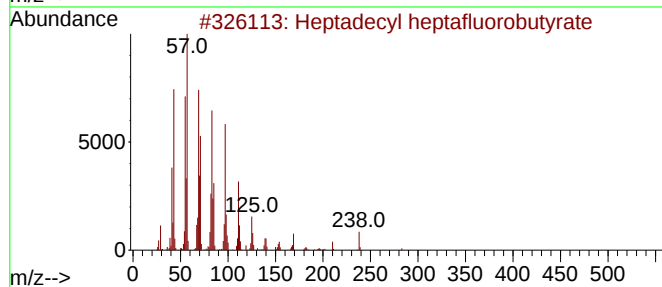
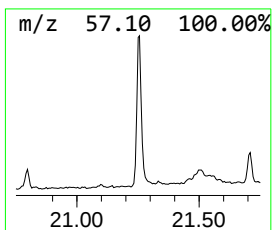
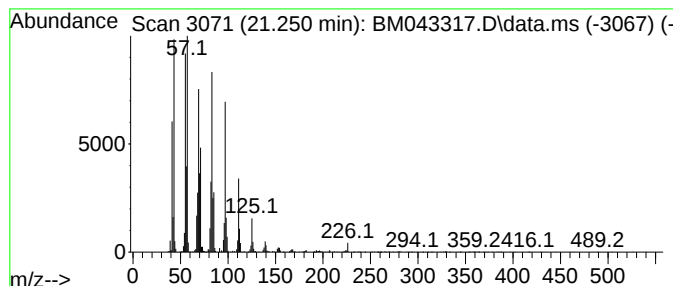
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.250	9.79 ng/ul	1869080	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3			Behenic alcohol	326	C22H46O	000661-19-8	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
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Sample : 05690-03
Misc :
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Instrument :
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ClientSampleId :
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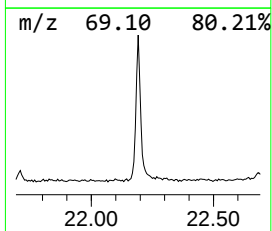
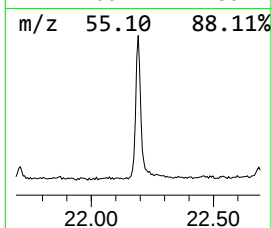
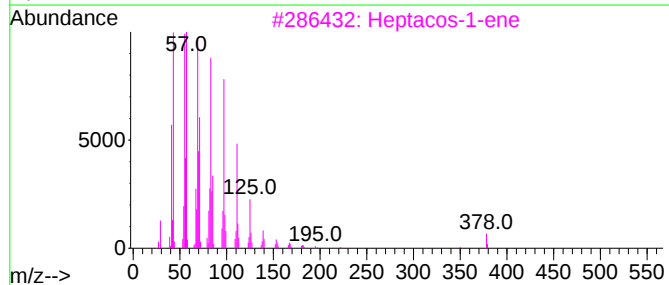
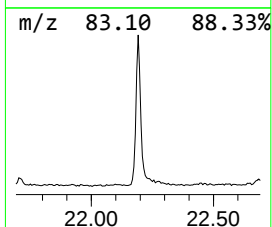
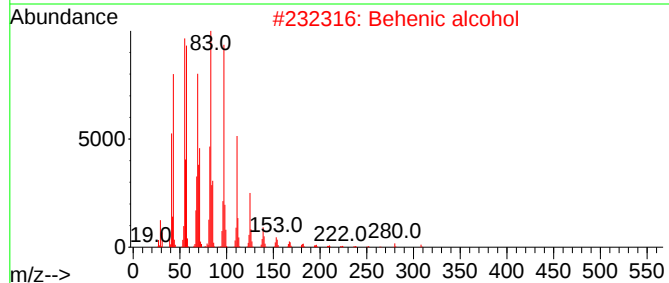
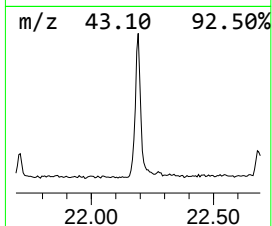
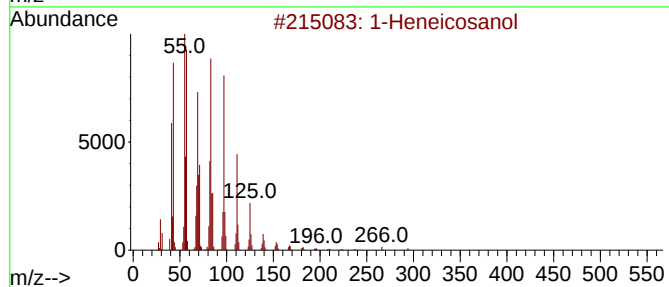
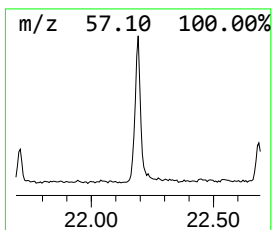
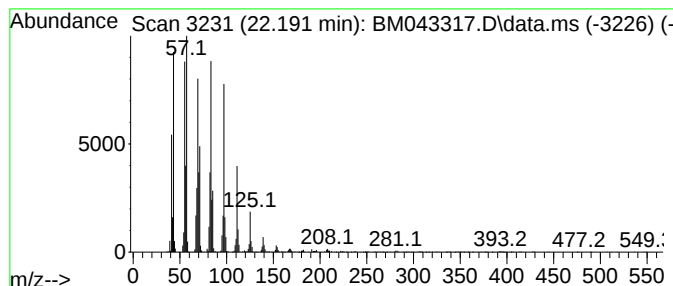
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.192	9.38 ng/ul	1791720	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Heptacos-1-ene	378	C27H54	015306-27-1	94
4			Nonacos-1-ene	406	C29H58	018835-35-3	94
5			n-Nonadecanol-1	284	C19H40O	001454-84-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
Operator : MA/JU
Sample : 05690-03
Misc :
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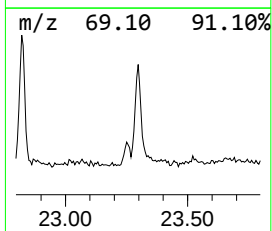
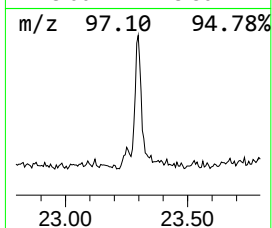
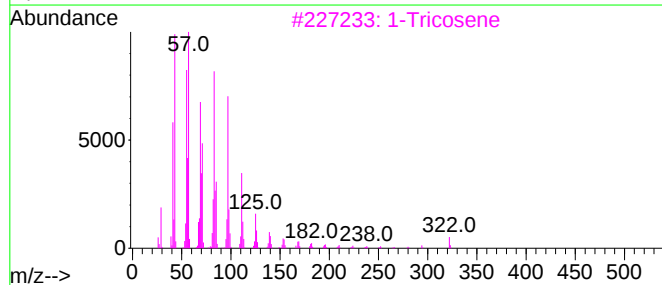
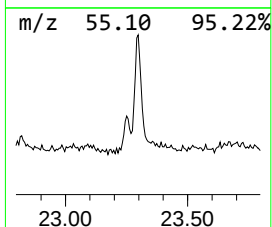
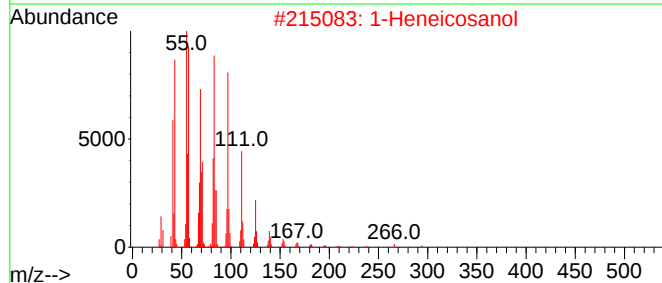
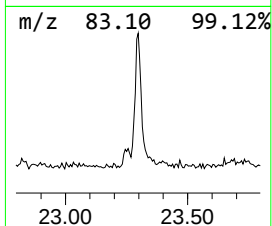
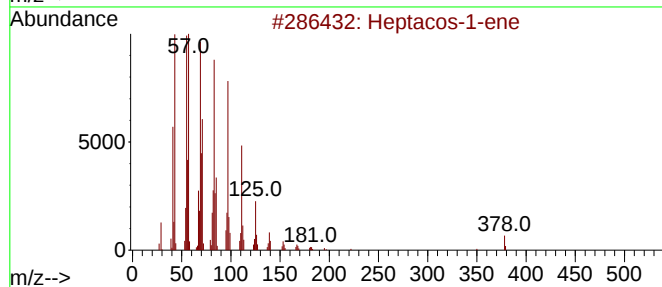
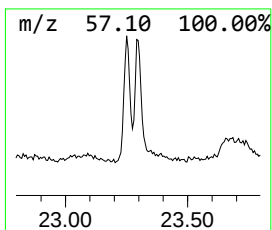
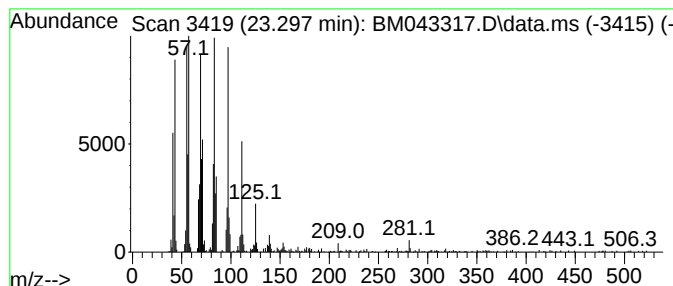
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.297	4.18 ng/ul	661369	Perylene-d12	23.956	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacos-1-ene	378	C27H54	015306-27-1	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	1-Tricosene	322	C23H46	018835-32-0	93
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	92
5	1-Hexacosanol	382	C26H54O	000506-52-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
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Instrument :
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ClientSampleId :
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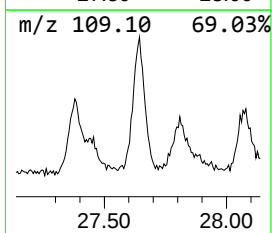
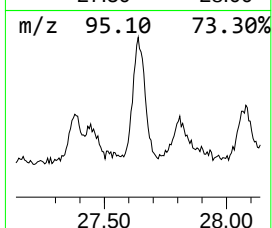
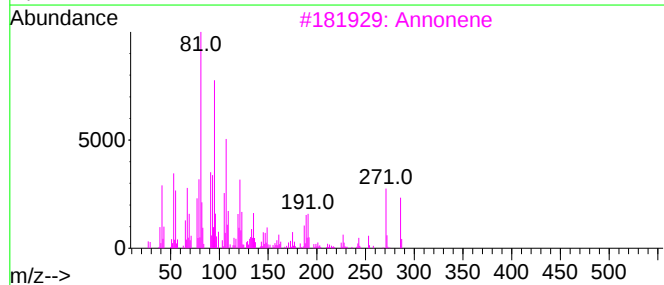
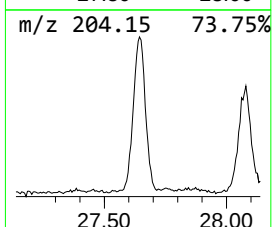
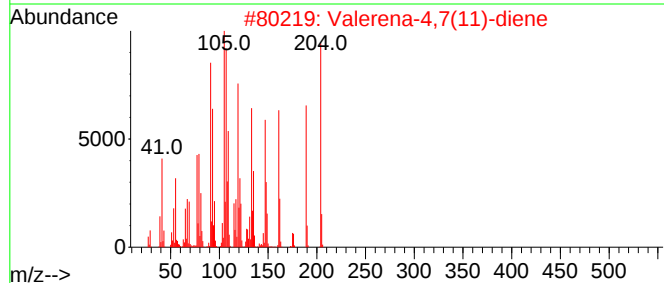
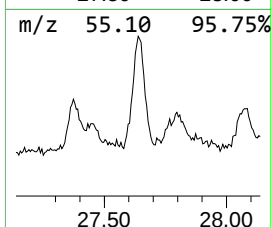
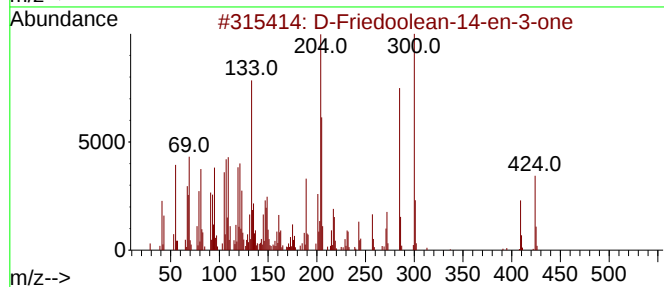
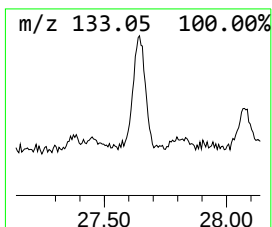
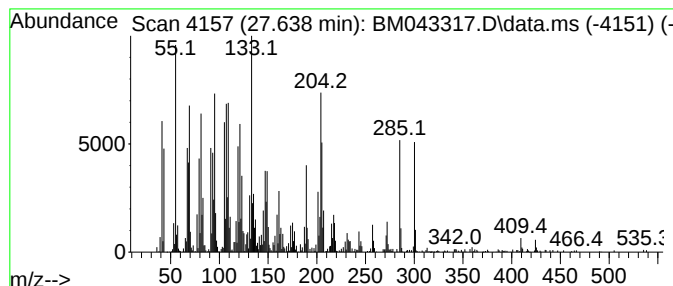
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 D-Friedoolean-14-en-3-one Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.638	13.60 ng/ul	2149680	Perylene-d12	23.956

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D-Friedoolean-14-en-3-one	424	C30H48O	000514-07-8	87
2			Valerena-4,7(11)-diene	204	C15H24	351222-66-7	64
3			Annonene	286	C20H30O	038611-50-6	47
4			Dimethyl(bis[(4,8,8-trimethyldec...	500	C32H56O2Si	1000364-69-0	43
5			Olean-18-ene	410	C30H50	000432-11-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121323\
Data File : BM043317.D
Acq On : 14 Dec 2023 00:37
Operator : MA/JU
Sample : 05690-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLH2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120723.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Aceto phenone	8.833	12.7	ng/ul	1516320	1	7.998	2389850	20.0
Benzoic acid	10.045	2.1	ng/ul	378322	2	10.804	3647240	20.0
Dibenzo furan	15.021	2.6	ng/ul	603642	3	14.627	4631670	20.0
n-Hexadecanoic ...	18.262	5.1	ng/ul	1232900	4	17.362	4871470	20.0
Heptadecyl hept...	21.250	9.8	ng/ul	1869080	5	21.533	3819530	20.0
1-Heneicosanol	22.192	9.4	ng/ul	1791720	5	21.533	3819530	20.0
(DEL) Alkane: S...	23.297	4.2	ng/ul	661369	6	23.956	3160750	20.0
D-Friedoolean-1...	27.638	13.6	ng/ul	2149680	6	23.956	3160750	20.0