

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
 Data File : BM034991.D
 Acq On : 11 May 2022 18:44
 Operator : CG/JU
 Sample : N2603-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW491

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-BM051022.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034991.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	44	48	54	rVB	71697	94478	1.56%	0.128%
2	3.646	92	95	102	rBV	16394	22938	0.38%	0.031%
3	3.781	114	118	131	rBV	420553	636685	10.52%	0.861%
4	3.887	132	136	142	rVV	89198	117938	1.95%	0.160%
5	4.022	155	159	164	rVB2	14606	22343	0.37%	0.030%
6	4.110	171	174	178	rVB	18781	24164	0.40%	0.033%
7	4.669	265	269	273	rVB3	13237	18695	0.31%	0.025%
8	5.028	325	330	335	rBV	52177	74044	1.22%	0.100%
9	5.122	343	346	353	rVB2	15484	22895	0.38%	0.031%
10	6.904	644	649	657	rBV4	14093	27380	0.45%	0.037%
11	7.081	672	679	690	rBV	1113177	1792690	29.62%	2.425%
12	7.251	701	708	717	rBV	877030	1354357	22.38%	1.832%
13	7.451	735	742	754	rBV	1203572	1904671	31.47%	2.577%
14	7.557	754	760	765	rVV6	13284	29441	0.49%	0.040%
15	7.922	815	822	829	rBV	874358	1364212	22.54%	1.845%
16	8.157	856	862	870	rBV4	18326	33951	0.56%	0.046%
17	8.628	934	942	956	rBV	1030385	1681865	27.79%	2.275%
18	9.075	1010	1018	1028	rBV	1075434	1789425	29.56%	2.421%
19	9.251	1041	1048	1057	rVB10	12503	23746	0.39%	0.032%
20	9.798	1131	1141	1157	rBV	901534	1531470	25.30%	2.072%
21	10.339	1226	1233	1247	rBV	1276843	2144783	35.43%	2.901%
22	10.504	1254	1261	1269	rBV	214135	344002	5.68%	0.465%
23	10.722	1290	1298	1304	rBV	1121001	1866513	30.84%	2.525%
24	10.775	1304	1307	1313	rVV	23179	36770	0.61%	0.050%
25	10.851	1313	1320	1338	rVV	605536	1123501	18.56%	1.520%
26	11.516	1427	1433	1439	rBV8	11373	23420	0.39%	0.032%
27	11.692	1458	1463	1471	rBV2	15853	29078	0.48%	0.039%
28	12.098	1526	1532	1538	rVV	24589	42597	0.70%	0.058%
29	12.310	1561	1568	1575	rVV	25498	46184	0.76%	0.062%
30	12.380	1575	1580	1587	rVB	25061	39645	0.65%	0.054%
31	12.598	1613	1617	1623	rVB	19671	30630	0.51%	0.041%
32	13.168	1709	1714	1719	rVB2	15707	25455	0.42%	0.034%
33	13.392	1748	1752	1756	rVB3	19551	25826	0.43%	0.035%
34	13.874	1829	1834	1839	rBV4	17184	27956	0.46%	0.038%
35	13.963	1839	1849	1855	rBV	2281586	3031748	50.09%	4.101%
36	14.245	1891	1897	1910	rVB	2532970	3618098	59.77%	4.894%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
 Title : SVOA CALIBRATION

37	14.463	1931	1934	1938	rBV3	14585	19370	0.32%	0.026%
38	14.551	1938	1949	1955	rBV	1503479	2266135	37.44%	3.066%
39	14.615	1955	1960	1968	rVB	115309	174432	2.88%	0.236%
40	14.739	1974	1981	1992	rBV	876392	1378715	22.78%	1.865%
41	14.892	2003	2007	2012	rBV4	20626	36547	0.60%	0.049%
42	14.951	2012	2017	2025	rVB	83401	156918	2.59%	0.212%
43	15.186	2054	2057	2066	rVB4	16544	38502	0.64%	0.052%
44	15.368	2081	2088	2092	rVV2	49388	85715	1.42%	0.116%
45	15.415	2092	2096	2100	rVB	44953	55310	0.91%	0.075%
46	15.474	2100	2106	2110	rBV7	10364	19413	0.32%	0.026%
47	15.545	2110	2118	2123	rBV	3186391	4500340	74.35%	6.088%
48	15.598	2123	2127	2131	rVB	150074	197683	3.27%	0.267%
49	15.657	2131	2137	2145	rBV	976529	1288499	21.29%	1.743%
50	15.780	2151	2158	2163	rBV3	38692	91952	1.52%	0.124%
51	15.898	2174	2178	2184	rVB3	31994	63952	1.06%	0.087%
52	15.968	2187	2190	2197	rVB5	17006	27735	0.46%	0.038%
53	16.039	2197	2202	2208	rBV	52851	97666	1.61%	0.132%
54	16.121	2214	2216	2221	rVB6	16255	21007	0.35%	0.028%
55	16.245	2230	2237	2245	rBV	2519324	3241604	53.55%	4.385%
56	16.445	2267	2271	2275	rVB5	15204	25519	0.42%	0.035%
57	16.604	2292	2298	2303	rVV5	27101	59184	0.98%	0.080%
58	16.668	2303	2309	2312	rVV4	25095	52035	0.86%	0.070%
59	16.733	2316	2320	2325	rBV	62885	88530	1.46%	0.120%
60	16.845	2334	2339	2342	rBV2	32374	50661	0.84%	0.069%
61	17.068	2372	2377	2381	rBV	235241	321101	5.30%	0.434%
62	17.115	2381	2385	2389	rVB2	80689	125998	2.08%	0.170%
63	17.221	2400	2403	2408	rVB2	43488	53081	0.88%	0.072%
64	17.292	2409	2415	2419	rBV	1691703	2438809	40.29%	3.299%
65	17.333	2419	2422	2427	rVB	1036274	1320173	21.81%	1.786%
66	17.392	2427	2432	2436	rBV2	3177720	4440135	73.35%	6.007%
67	17.486	2444	2448	2454	rVB4	49310	89207	1.47%	0.121%
68	17.539	2454	2457	2460	rBV3	23619	28849	0.48%	0.039%
69	17.586	2462	2465	2466	rBV2	19790	22981	0.38%	0.031%
70	17.692	2476	2483	2487	rBV2	89331	146306	2.42%	0.198%
71	17.803	2498	2502	2511	rBV2	126760	181655	3.00%	0.246%
72	17.974	2528	2531	2535	rVB3	46666	61614	1.02%	0.083%
73	18.080	2546	2549	2554	rBV4	21287	35737	0.59%	0.048%
74	18.133	2555	2558	2561	rVV3	32449	39479	0.65%	0.053%
75	18.198	2561	2569	2576	rVV2	721975	1175095	19.41%	1.590%
76	18.262	2578	2580	2584	rVV	50974	59297	0.98%	0.080%

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77	18.298	2584	2586	2592	rVB2	48916	73385	1.21%	0.099%
78	18.368	2592	2598	2602	rBV2	150244	253985	4.20%	0.344%
79	18.498	2616	2620	2627	rBV4	58940	110234	1.82%	0.149%
80	18.627	2639	2642	2645	rVB2	32098	33716	0.56%	0.046%
81	18.668	2645	2649	2653	rVV	77223	105164	1.74%	0.142%
82	18.703	2653	2655	2659	rVB2	33644	38948	0.64%	0.053%
83	18.856	2679	2681	2686	rBV6	19090	22404	0.37%	0.030%
84	18.962	2697	2699	2703	rVB3	24750	28011	0.46%	0.038%
85	19.080	2715	2719	2724	rBV3	71441	97703	1.61%	0.132%
86	19.245	2745	2747	2751	rVB2	48018	54547	0.90%	0.074%
87	19.345	2756	2764	2774	rVV	1357683	2018561	33.35%	2.731%
88	19.468	2781	2785	2794	rVB	318245	412703	6.82%	0.558%
89	19.680	2815	2821	2825	rBV	4245258	6052988	100.00%	8.188%
90	19.886	2853	2856	2862	rVB	60163	77296	1.28%	0.105%
91	20.203	2906	2910	2913	rBV2	184587	209979	3.47%	0.284%
92	20.303	2923	2927	2931	rBV	171995	229852	3.80%	0.311%
93	20.609	2976	2979	2984	rVB2	96264	113823	1.88%	0.154%
94	21.186	3074	3077	3084	rVB3	964454	1297832	21.44%	1.756%
95	21.468	3119	3125	3129	rBV2	2121196	3261485	53.88%	4.412%
96	21.503	3129	3131	3140	rVB	485326	583043	9.63%	0.789%
97	22.103	3229	3233	3241	rBV3	144300	269658	4.45%	0.365%
98	23.138	3403	3409	3413	rBV3	429457	1020825	16.86%	1.381%
99	23.703	3498	3505	3511	rBV2	2755682	5256918	86.85%	7.111%
100	23.856	3524	3531	3538	rBV2	1344290	2698376	44.58%	3.650%

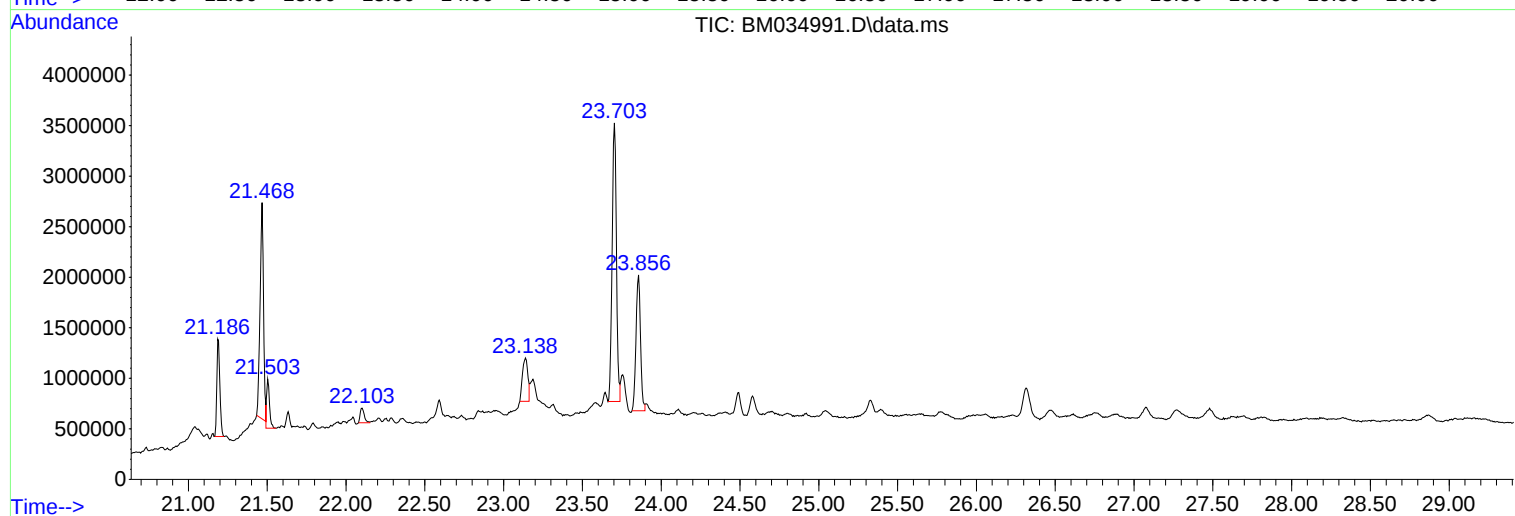
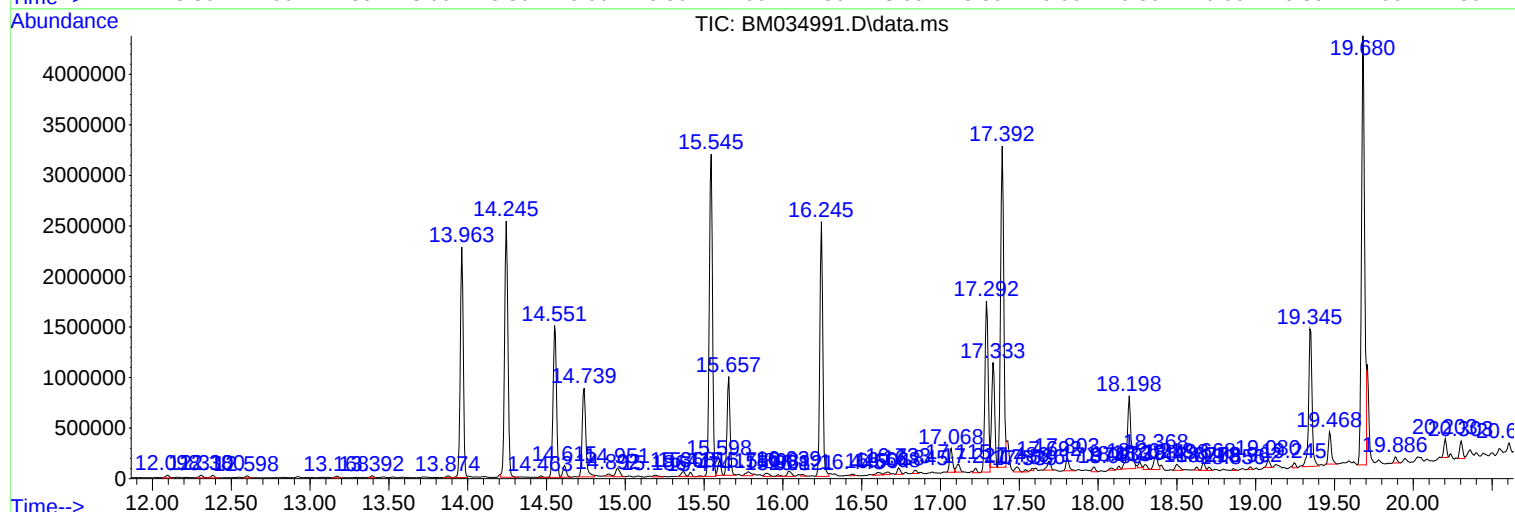
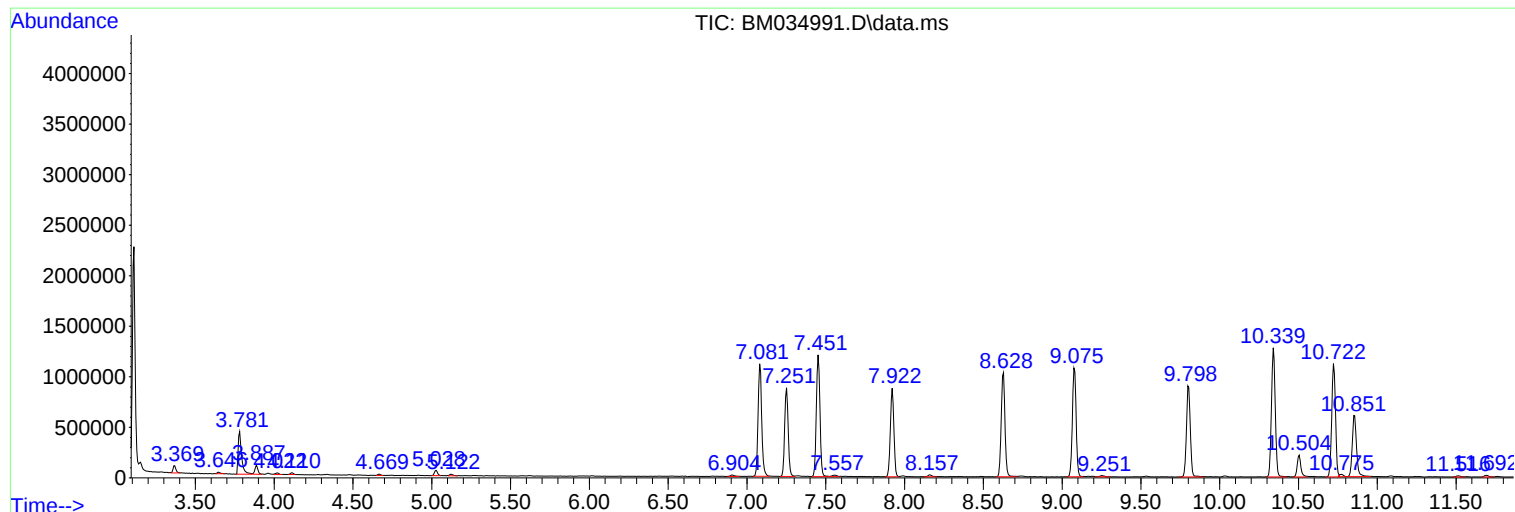
Sum of corrected areas: 73921971

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

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



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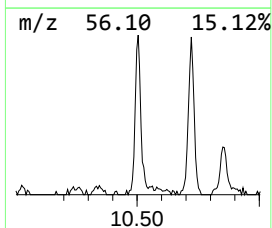
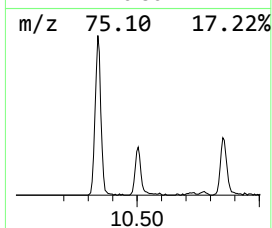
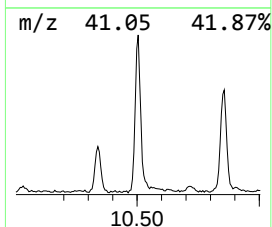
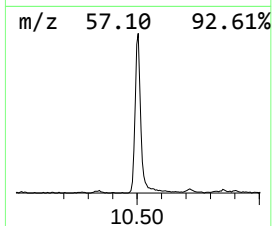
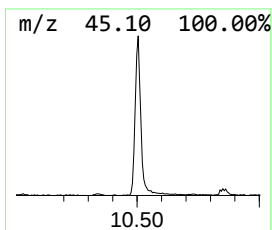
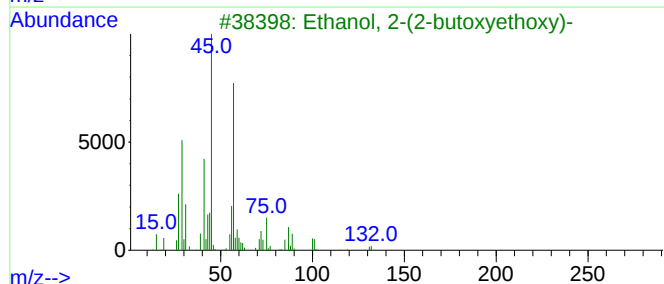
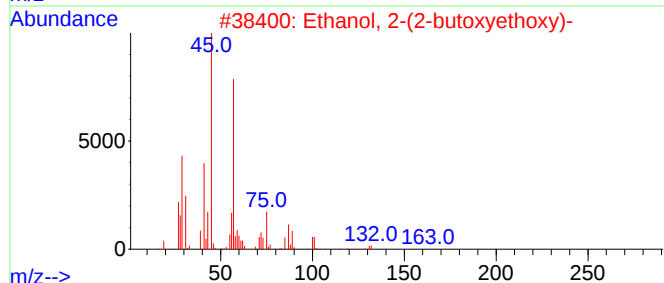
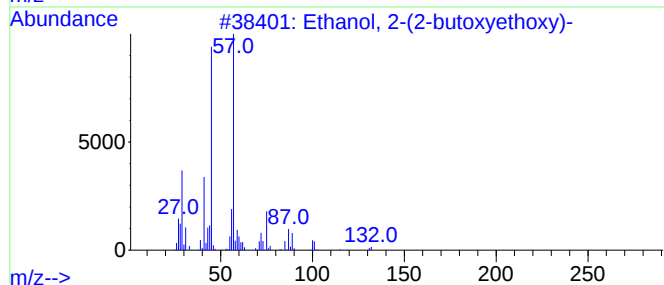
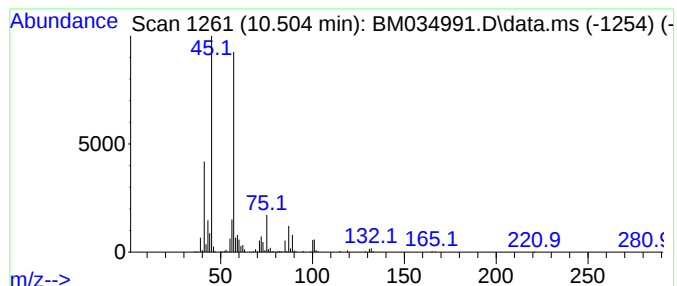
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.504	3.69 ng/ul	344002	Naphthalene-d8	10.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
4			Di-sec-Butyl ether	130	C8H18O	006863-58-7	53
5			Butyl lactate	146	C7H14O3	000138-22-7	50



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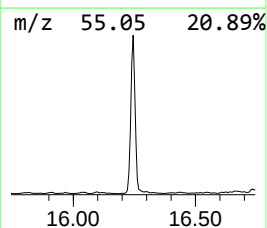
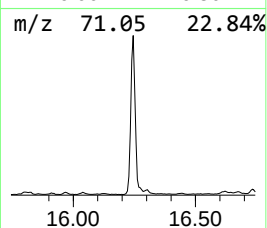
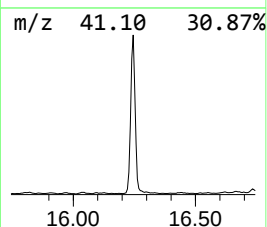
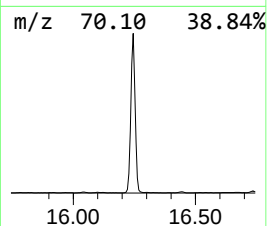
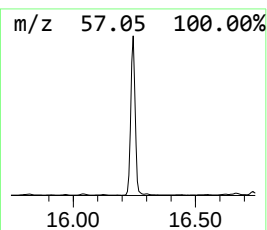
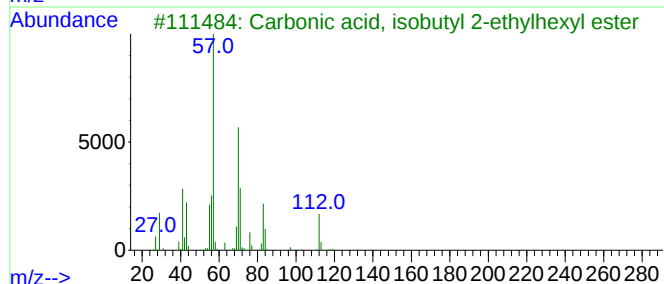
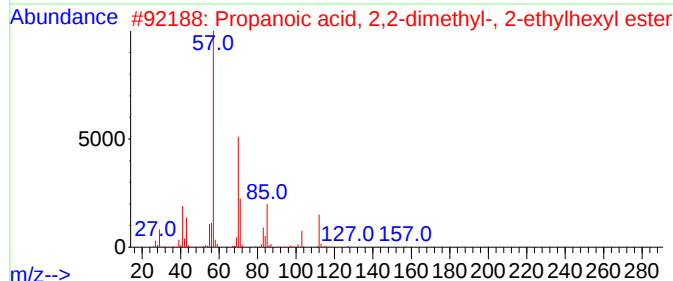
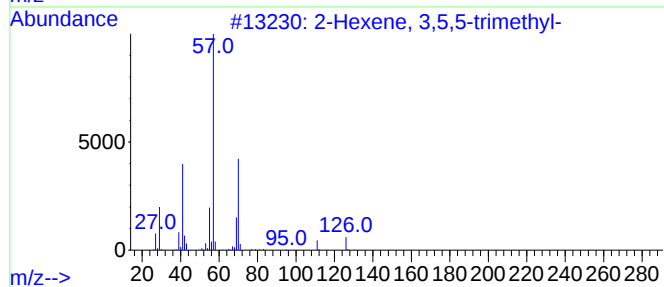
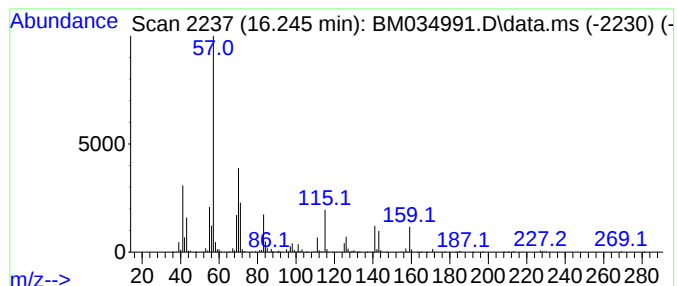
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.245	26.58 ng/ul	3241600	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 3,5,5-trimethyl-		126	C9H18		026456-76-8	43
2	Propanoic acid, 2,2-dimethyl-, 2...		214	C13H26O2		016387-18-1	37
3	Carbonic acid, isobutyl 2-ethylh...		230	C13H26O3		1000383-13-3	37
4	Carbonic acid, allyl 2-ethylhexy...		214	C12H22O3		1000357-80-6	35
5	Hexanoic acid, 3,5,5-trimethyl-,...		368	C24H48O2		1000406-27-5	35



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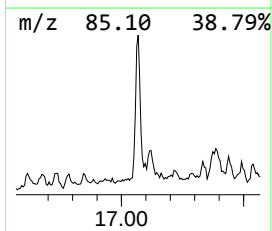
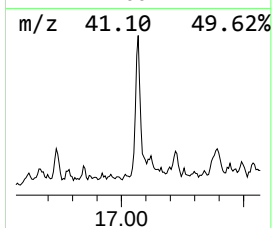
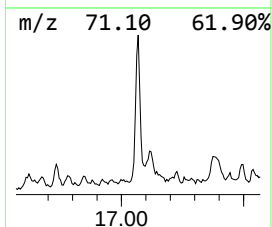
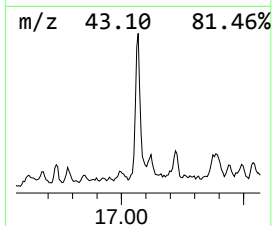
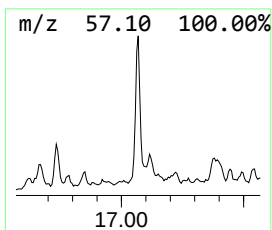
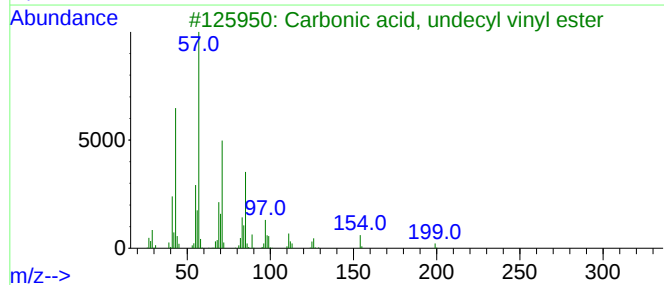
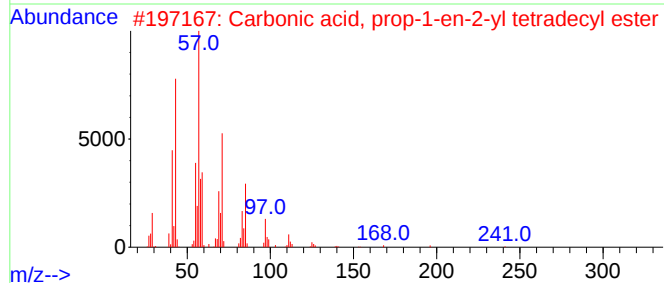
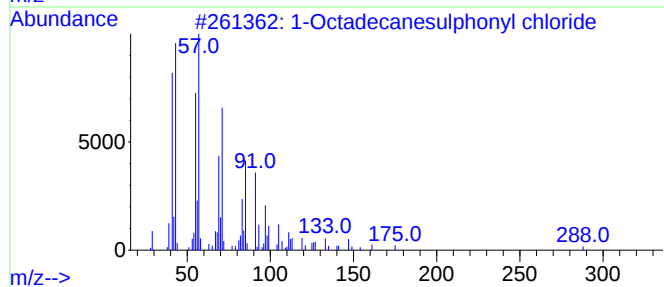
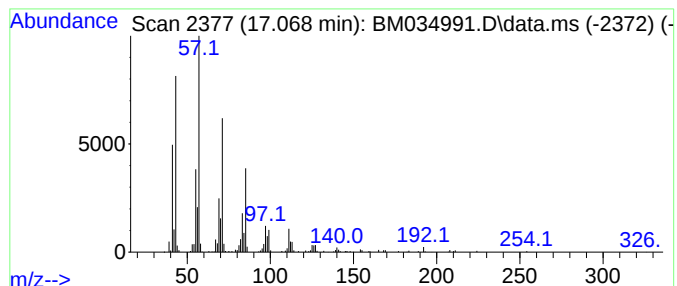
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EW491

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Octadecanesulphonyl chloride Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.068	2.63 ng/ul	321101	Phenanthrene-d10	17.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	90
2	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	87
3	Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	86
4	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	86
5	Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034991.D
Acq On : 11 May 2022 18:44
Operator : CG/JU
Sample : N2603-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW491

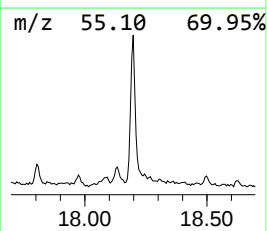
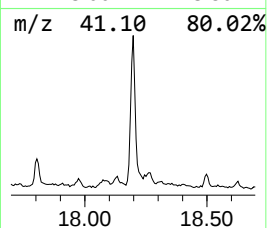
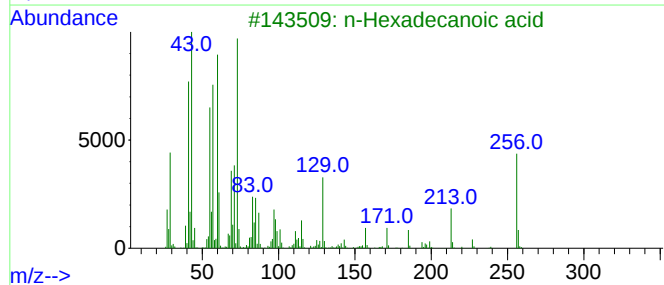
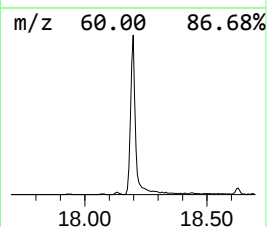
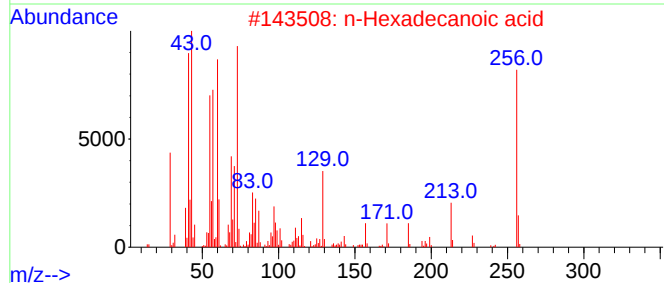
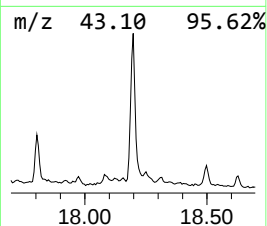
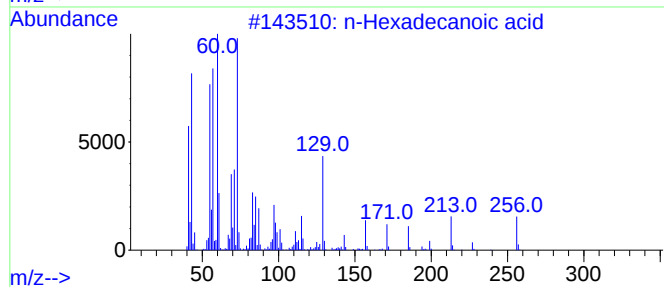
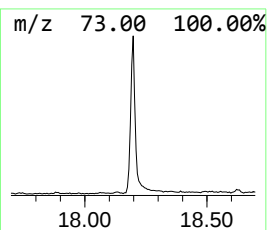
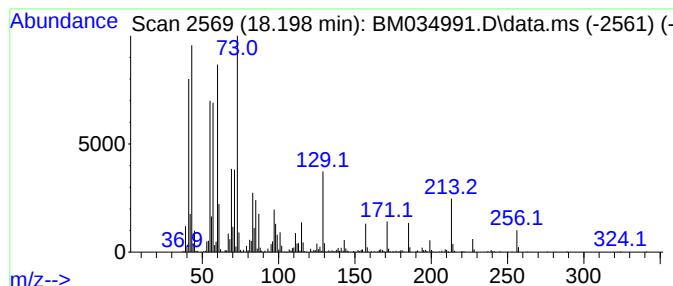
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.198	9.64 ng/ul	1175100	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034991.D
Acq On : 11 May 2022 18:44
Operator : CG/JU
Sample : N2603-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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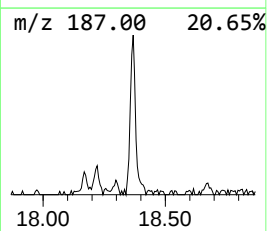
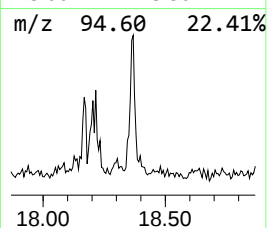
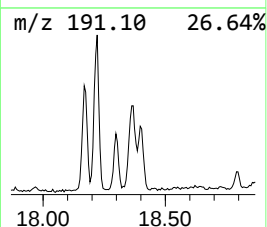
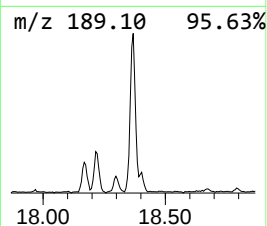
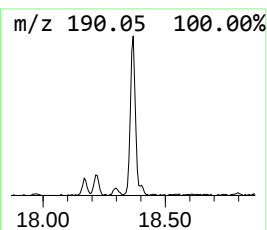
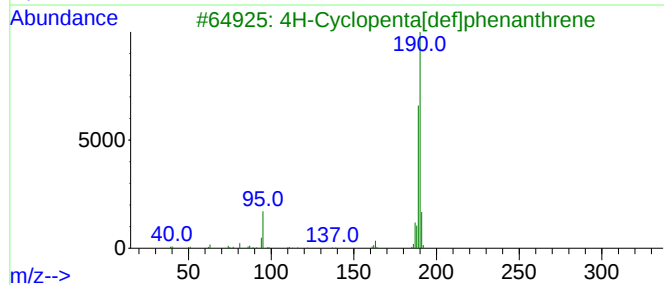
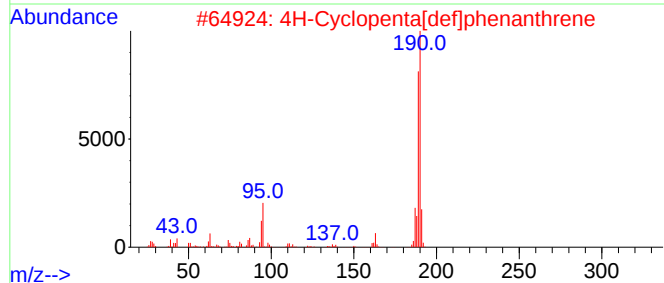
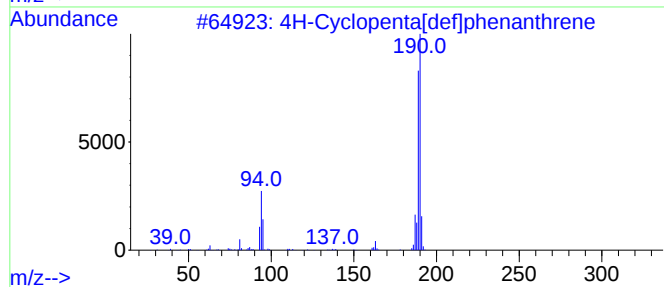
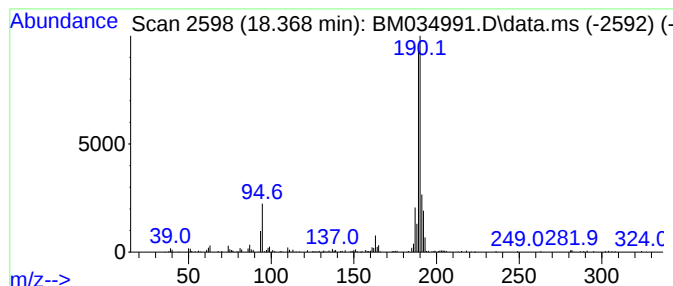
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.368	2.08 ng/ul	253985	Phenanthrene-d10	17.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034991.D
Acq On : 11 May 2022 18:44
Operator : CG/JU
Sample : N2603-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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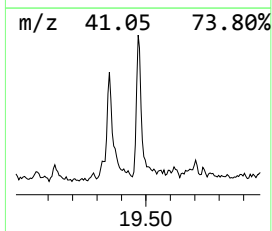
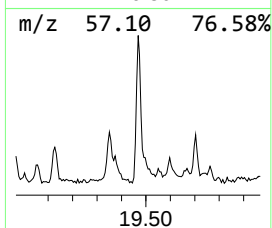
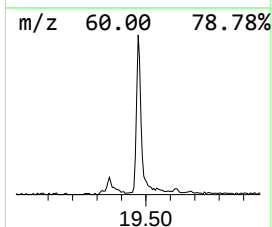
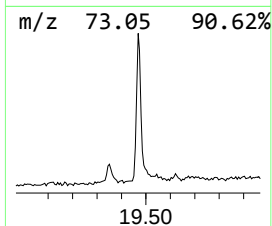
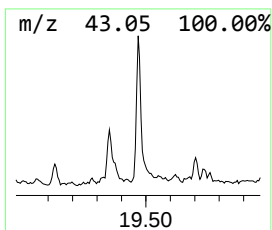
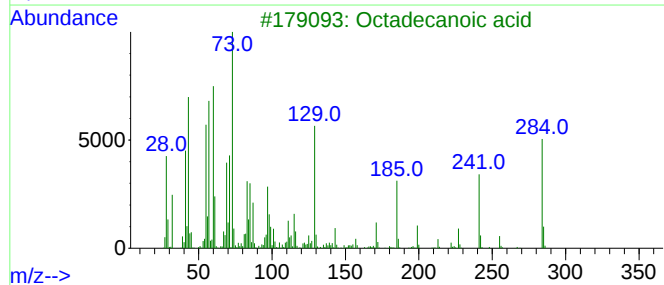
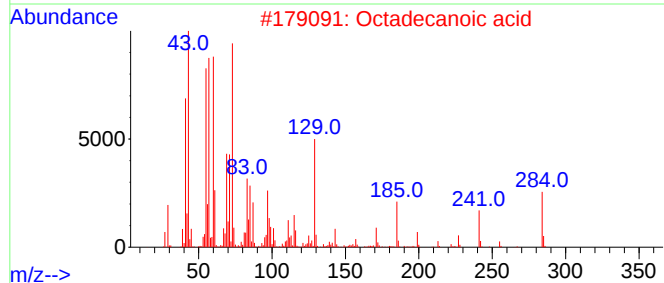
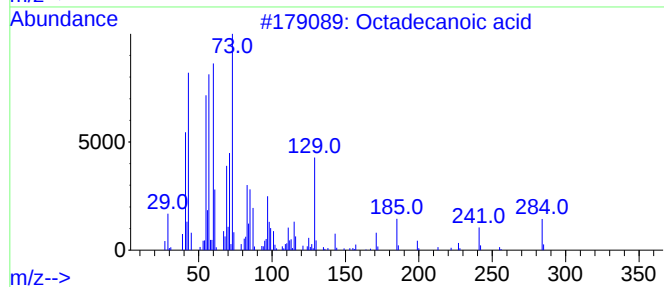
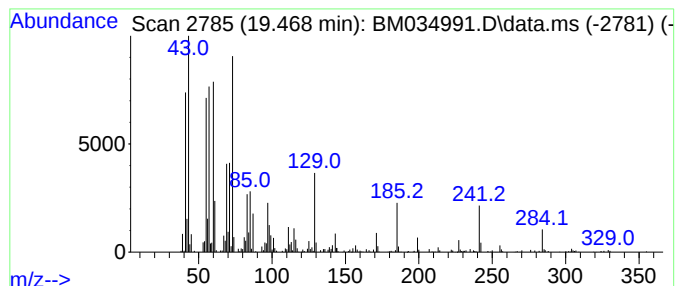
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.468	2.53 ng/ul	412703	Chrysene-d12	21.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034991.D
Acq On : 11 May 2022 18:44
Operator : CG/JU
Sample : N2603-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW491

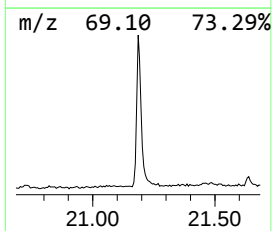
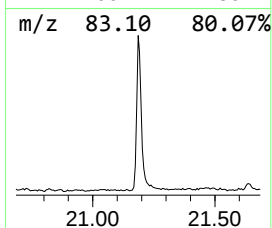
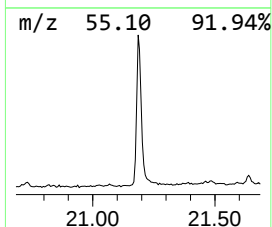
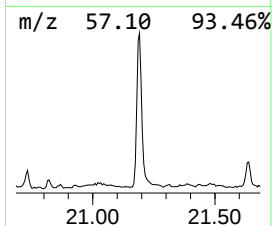
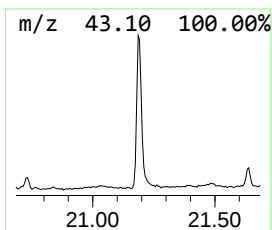
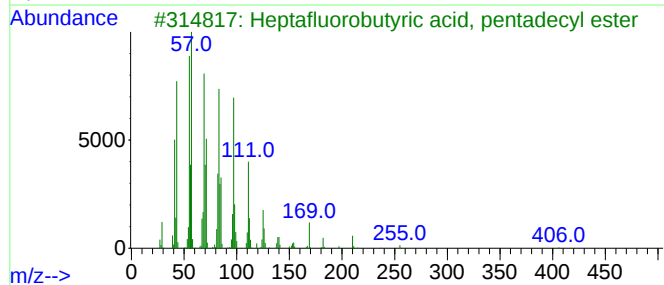
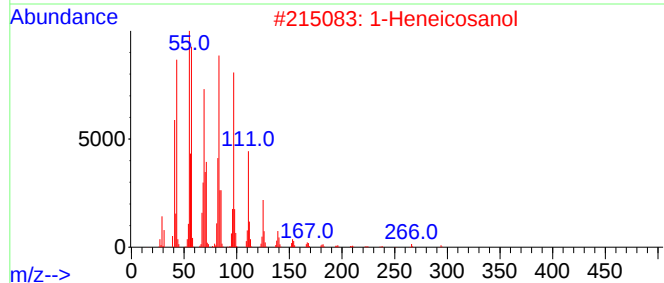
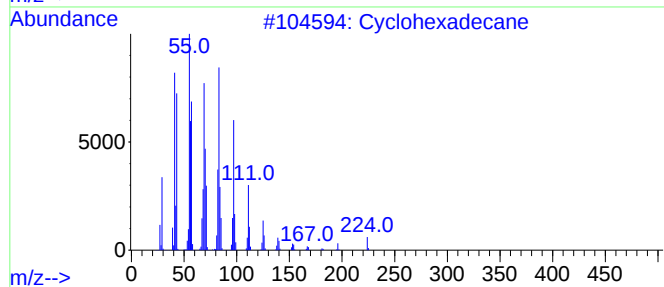
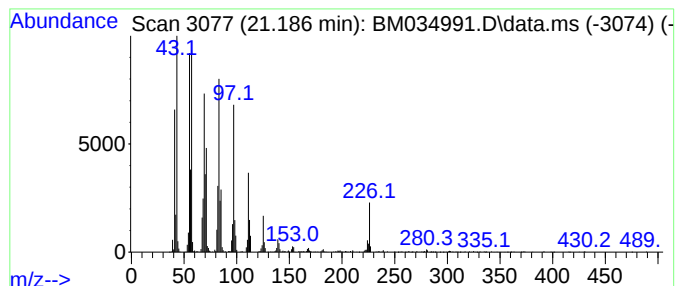
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.186 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.186	7.96 ng/ul	1297830	Chrysene-d12	21.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	93
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
4			Pentacos-1-ene	350	C25H50	016980-85-1	93
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034991.D
Acq On : 11 May 2022 18:44
Operator : CG/JU
Sample : N2603-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.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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(DEL) Alkane: S...	16.245	26.6	ng/ul	3241600	4	17.292	2438810	20.0
1-Octadecanesul...	17.068	2.6	ng/ul	321101	4	17.292	2438810	20.0
n-Hexadecanoic ...	18.198	9.6	ng/ul	1175100	4	17.292	2438810	20.0
4H-Cyclopenta[d...	18.368	2.1	ng/ul	253985	4	17.292	2438810	20.0
Octadecanoic acid	19.468	2.5	ng/ul	412703	5	21.468	3261490	20.0
(DEL) Alkane: C...	21.186	8.0	ng/ul	1297830	5	21.468	3261490	20.0