

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
 Data File : BP015187.D
 Acq On : 19 May 2023 23:11
 Operator : CG/JU
 Sample : 02689-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREM3

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_P\Methods\SFAM-EPA-BP051123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015187.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	38	42	52	rVB	50871	76253	1.04%	0.094%
2	3.875	114	117	132	rBV	450228	680894	9.30%	0.837%
3	3.987	132	136	139	rVB	35726	46742	0.64%	0.057%
4	4.110	152	157	163	rVB	37269	55735	0.76%	0.068%
5	4.210	170	174	181	rVB2	14257	20078	0.27%	0.025%
6	4.599	236	240	246	rBV5	15164	26860	0.37%	0.033%
7	4.669	248	252	259	rVB	68488	101686	1.39%	0.125%
8	5.163	332	336	345	rVB3	30485	61539	0.84%	0.076%
9	5.287	350	357	361	rVB7	9873	19172	0.26%	0.024%
10	5.428	376	381	387	rBV5	8870	19572	0.27%	0.024%
11	5.657	416	420	425	rBV	17168	22998	0.31%	0.028%
12	5.781	436	441	450	rVB	32395	52122	0.71%	0.064%
13	6.781	607	611	617	rVB7	8648	13486	0.18%	0.017%
14	7.228	682	687	690	rBV	57458	86923	1.19%	0.107%
15	7.275	690	695	714	rVV	1095512	1821685	24.89%	2.238%
16	7.446	714	724	736	rVV	981915	1612546	22.03%	1.981%
17	7.651	752	759	768	rBV	1172244	1840734	25.15%	2.261%
18	7.728	770	772	780	rVB6	9059	17879	0.24%	0.022%
19	7.846	788	792	799	rVB2	8847	17158	0.23%	0.021%
20	7.940	802	808	814	rVB2	9261	16978	0.23%	0.021%
21	8.128	831	840	849	rBV	945421	1533172	20.95%	1.884%
22	8.340	873	876	880	rBV5	9594	12907	0.18%	0.016%
23	8.834	952	960	975	rBV	1223035	2086560	28.51%	2.564%
24	8.963	977	982	989	rVB2	17247	31335	0.43%	0.038%
25	9.298	1031	1039	1051	rBV	1113878	1917563	26.20%	2.356%
26	9.410	1054	1058	1062	rVV6	7961	13459	0.18%	0.017%
27	9.457	1062	1066	1072	rVB9	7170	13666	0.19%	0.017%
28	9.693	1098	1106	1114	rBV	178113	266700	3.64%	0.328%
29	10.028	1154	1163	1181	rBV	988316	1679354	22.94%	2.063%
30	10.569	1247	1255	1267	rBV	1474492	2512082	34.32%	3.086%
31	10.645	1267	1268	1277	rVB9	8287	16050	0.22%	0.020%
32	10.957	1312	1321	1327	rBV	1390830	2328190	31.81%	2.860%
33	11.004	1327	1329	1335	rVB5	7651	11496	0.16%	0.014%
34	11.092	1335	1344	1359	rBV	1205772	2159605	29.50%	2.653%
35	11.487	1405	1411	1418	rBV9	12749	26190	0.36%	0.032%
36	11.734	1448	1453	1457	rBV3	14693	30004	0.41%	0.037%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
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37	12.187	1524	1530	1538	rVB	125180	180341	2.46%	0.222%
38	12.345	1552	1557	1564	rVB	10773	18697	0.26%	0.023%
39	12.428	1566	1571	1578	rBV3	13178	22186	0.30%	0.027%
40	12.534	1582	1589	1595	rVV	22937	37855	0.52%	0.047%

41	12.604	1599	1601	1607	rVV6	7496	13096	0.18%	0.016%
42	12.751	1618	1626	1632	rBV	11517	20572	0.28%	0.025%
43	13.304	1711	1720	1728	rVB	35520	62487	0.85%	0.077%
44	13.575	1756	1766	1773	rVV8	13109	31906	0.44%	0.039%
45	13.645	1773	1778	1783	rVV	57815	92657	1.27%	0.114%

46	13.734	1787	1793	1799	rVB8	8684	23273	0.32%	0.029%
47	13.910	1815	1823	1834	rBV	93768	266176	3.64%	0.327%
48	14.086	1848	1853	1859	rVB7	10744	17676	0.24%	0.022%
49	14.169	1859	1867	1880	rBV	3006578	4291127	58.62%	5.272%
50	14.404	1902	1907	1909	rBV6	8527	12863	0.18%	0.016%

51	14.457	1909	1916	1928	rVV2	3195431	4862082	66.42%	5.973%
52	14.763	1961	1968	1976	rBV	2160463	3172769	43.34%	3.898%
53	14.822	1976	1978	1983	rVB6	9444	12398	0.17%	0.015%
54	14.957	1994	2001	2013	rBV	1084104	1759434	24.04%	2.162%
55	15.045	2013	2016	2023	rVV4	27537	58694	0.80%	0.072%

56	15.110	2023	2027	2031	rVV4	23013	42955	0.59%	0.053%
57	15.169	2032	2037	2046	rVB3	66065	119223	1.63%	0.146%
58	15.269	2049	2054	2059	rVB	32590	49627	0.68%	0.061%
59	15.486	2087	2091	2096	rBV8	10372	16876	0.23%	0.021%
60	15.710	2123	2129	2131	rBV	112516	149388	2.04%	0.184%

61	15.757	2131	2137	2143	rVB	4174909	6065532	82.86%	7.452%
62	15.875	2151	2157	2165	rBV	1168231	1644856	22.47%	2.021%
63	16.080	2188	2192	2193	rBV3	11683	13602	0.19%	0.017%
64	16.116	2193	2198	2204	rVB	479434	628897	8.59%	0.773%
65	16.304	2226	2230	2233	rBV6	11466	19434	0.27%	0.024%

66	16.445	2251	2254	2260	rBV7	11307	18162	0.25%	0.022%
67	16.680	2290	2294	2301	rVB2	42259	57675	0.79%	0.071%
68	16.892	2326	2330	2337	rBV3	22040	27699	0.38%	0.034%
69	16.957	2337	2341	2347	rVV	100143	128800	1.76%	0.158%
70	17.045	2352	2356	2360	rVB3	18647	24366	0.33%	0.030%

71	17.163	2373	2376	2379	rBV5	13687	18817	0.26%	0.023%
72	17.216	2380	2385	2389	rVV	267076	357817	4.89%	0.440%
73	17.257	2389	2392	2396	rVV3	28175	39817	0.54%	0.049%
74	17.310	2396	2401	2405	rVV3	26185	37242	0.51%	0.046%
75	17.392	2412	2415	2422	rVB8	13560	22082	0.30%	0.027%

76	17.516	2430	2436	2442	rVV	2660652	3722473	50.85%	4.573%
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77	17.616	2446	2453	2463	rVV	4671924	6625375	90.51%	8.140%
78	17.698	2463	2467	2476	rVB2	99290	155069	2.12%	0.191%
79	18.027	2519	2523	2528	rVB	101468	124074	1.70%	0.152%
80	18.116	2534	2538	2542	rBV3	46578	62029	0.85%	0.076%
81	18.257	2559	2562	2569	rBV9	22080	33591	0.46%	0.041%
82	18.374	2576	2582	2591	rBV	871908	1261381	17.23%	1.550%
83	18.745	2642	2645	2649	rBV	28629	30991	0.42%	0.038%
84	18.933	2673	2677	2681	rBV	87971	104762	1.43%	0.129%
85	19.251	2728	2731	2736	rBV	69204	87443	1.19%	0.107%
86	19.410	2755	2758	2762	rVB	70031	84223	1.15%	0.103%
87	19.480	2764	2770	2772	rBV	80267	142809	1.95%	0.175%
88	19.592	2785	2789	2801	rBV	279159	468908	6.41%	0.576%
89	19.692	2803	2806	2810	rVV	103398	129505	1.77%	0.159%
90	19.833	2824	2830	2839	rBV	5656052	7319965	100.00%	8.993%
91	20.368	2918	2921	2924	rVB	102617	106244	1.45%	0.131%
92	20.810	2993	2996	3000	rBV	104790	108848	1.49%	0.134%
93	20.974	3022	3024	3029	rBV	102105	116111	1.59%	0.143%
94	21.268	3070	3074	3081	rBV	930647	1360824	18.59%	1.672%
95	21.592	3124	3129	3139	rVB	2062673	2908297	39.73%	3.573%
96	22.198	3229	3232	3236	rBV	334615	409333	5.59%	0.503%
97	23.321	3419	3423	3430	rVB	521416	832750	11.38%	1.023%
98	23.986	3529	3536	3546	rVB2	2638184	5727241	78.24%	7.036%
99	24.151	3558	3564	3573	rVB	1302638	2737517	37.40%	3.363%
100	24.792	3668	3673	3682	rVB	476814	1046353	14.29%	1.286%

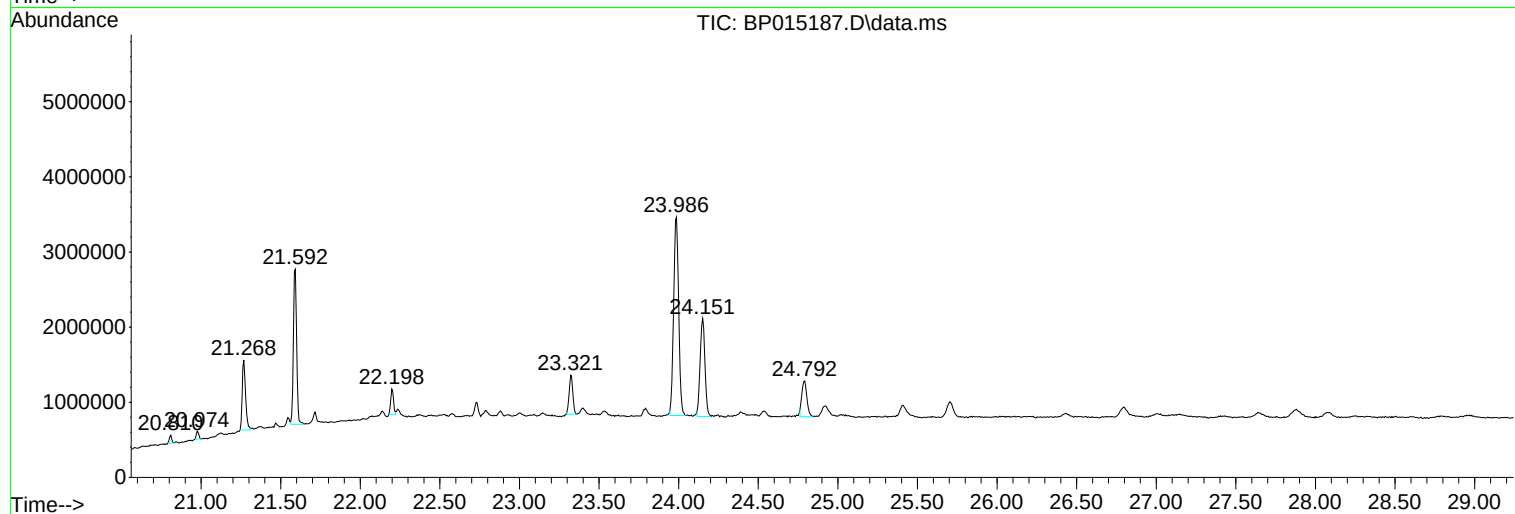
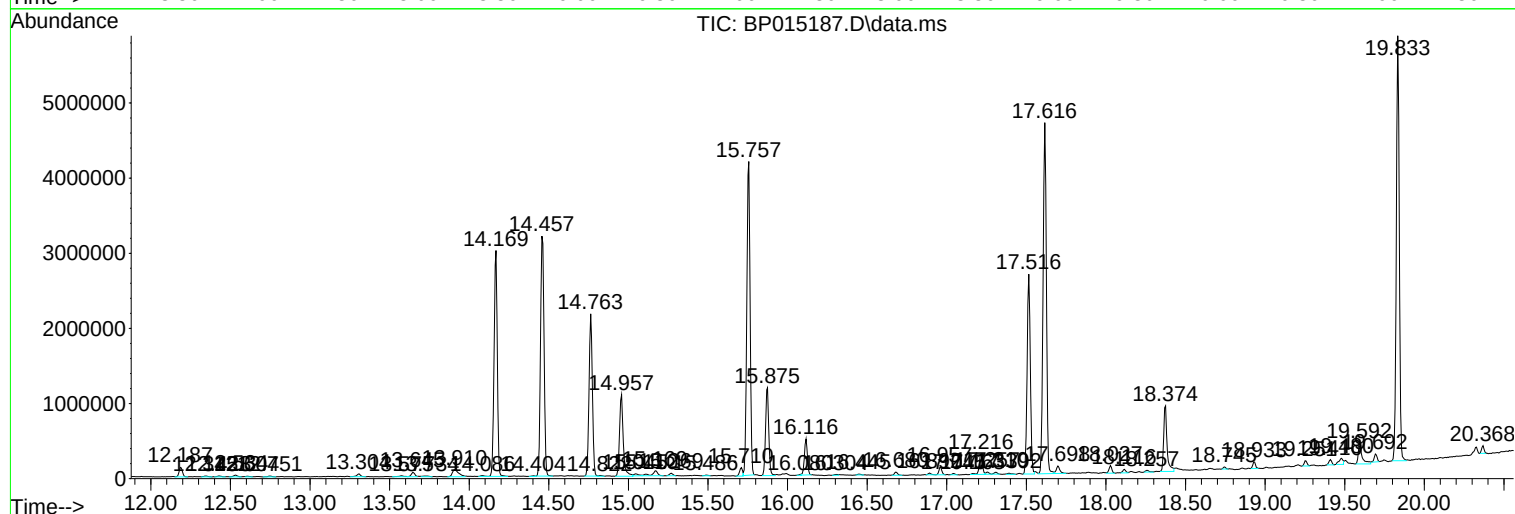
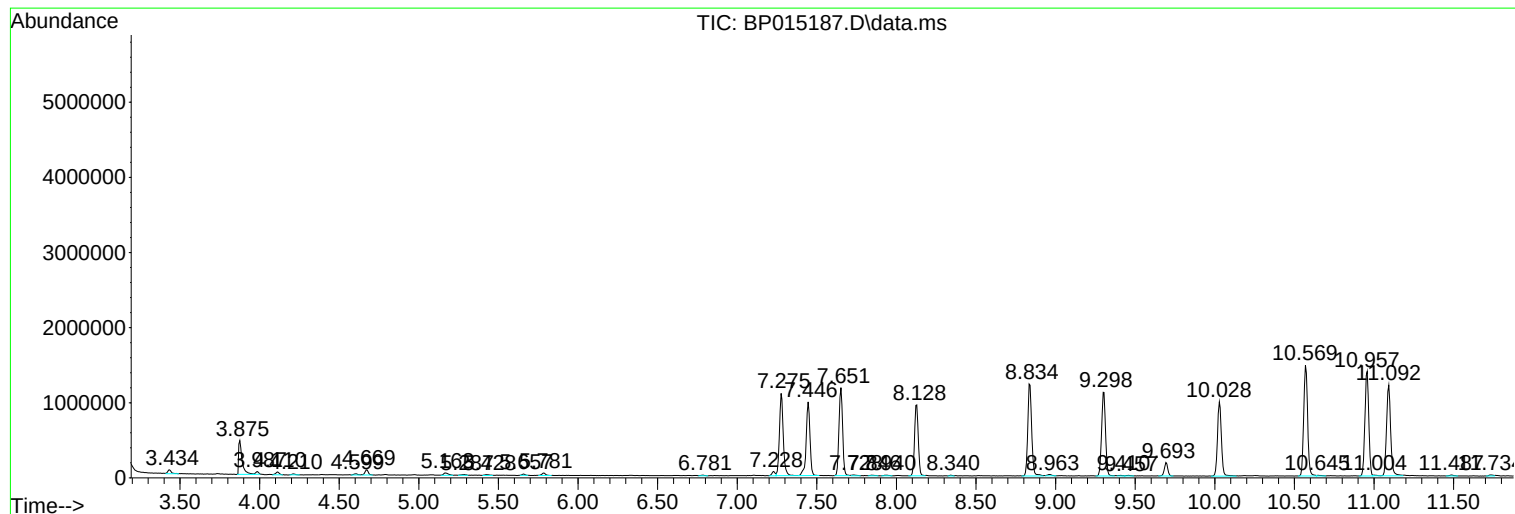
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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



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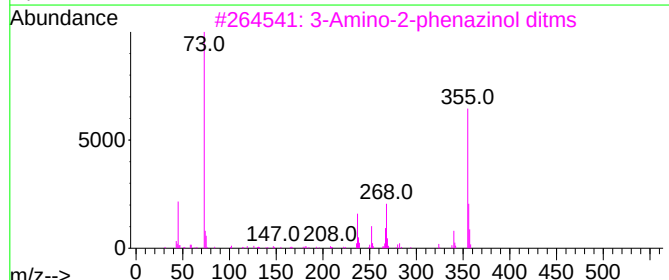
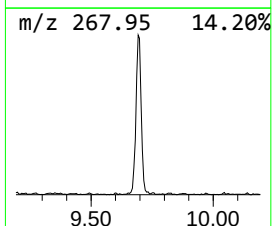
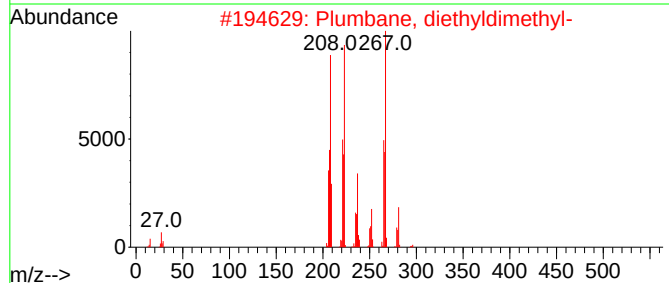
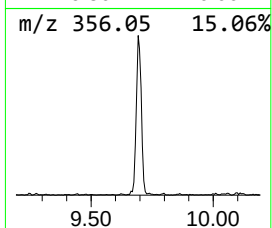
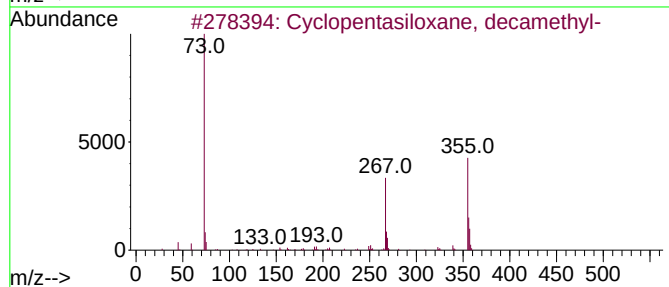
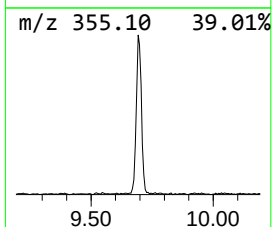
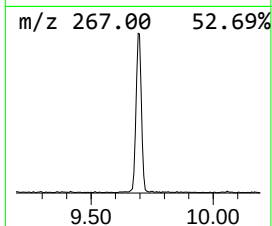
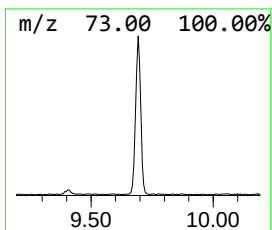
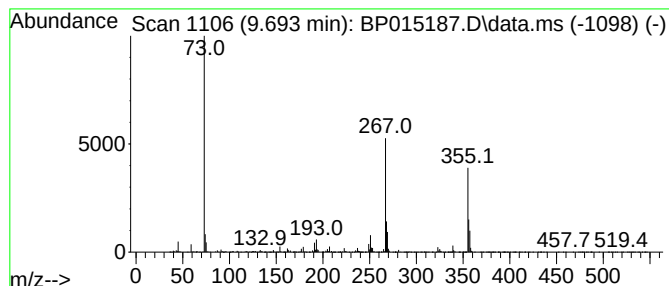
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopentasiloxane, decamet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.693	2.29 ng/ul	266700	Naphthalene-d8	10.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2			Plumbane, diethyldimethyl-	296	C6H16Pb	001762-27-2	91
3			3-Amino-2-phenazinol ditms	355	C18H25N3OSi2	1000332-15-5	47
4			4-(4-Fluorophenyl)-2,3-dihydro-1...	267	C12H14FN0SSi	1010501-61-0	43
5			Bamethan, 2TMS derivative	353	C18H35N02Si2	040629-66-1	43



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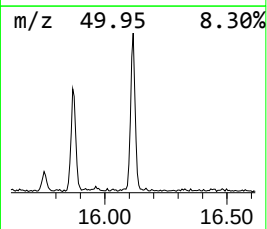
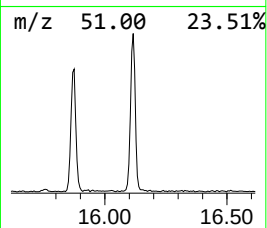
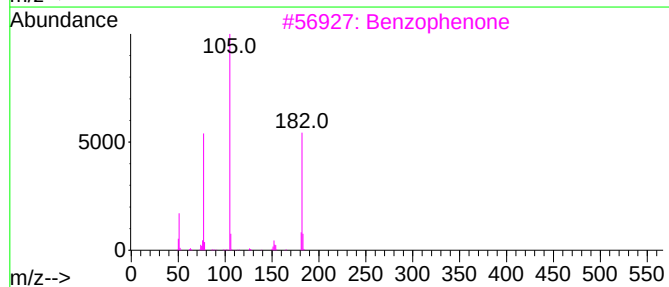
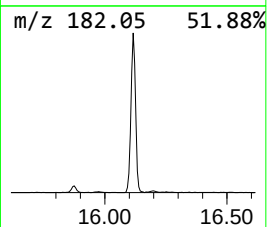
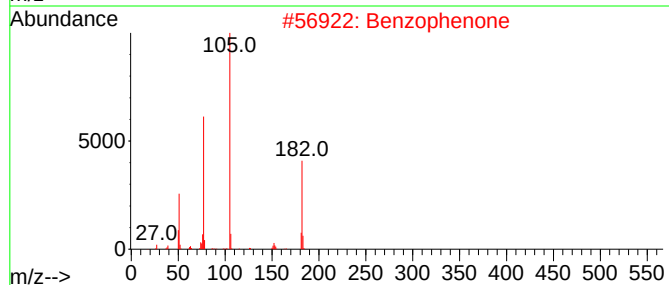
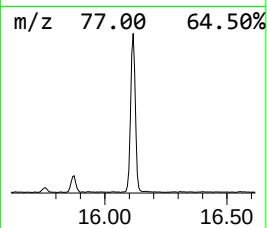
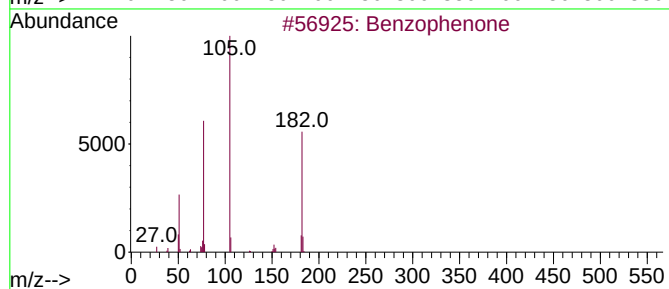
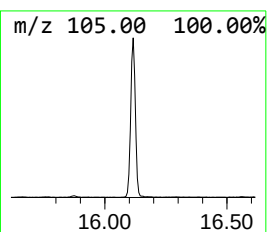
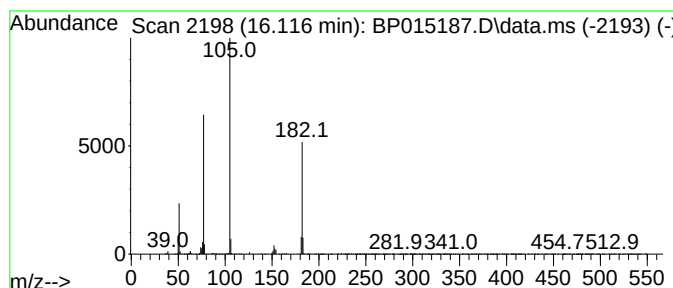
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.116	3.96 ng/ul	628897	Acenaphthene-d10	14.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	87
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64



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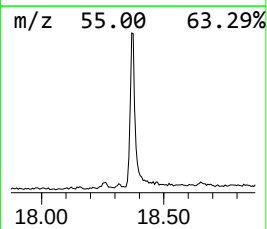
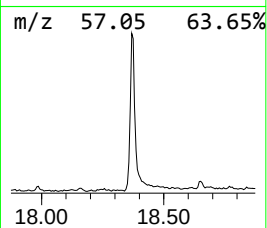
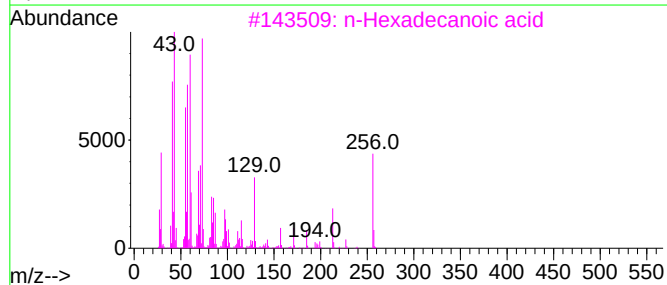
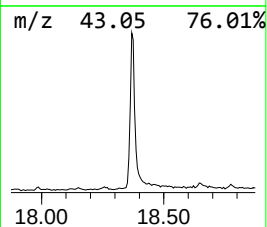
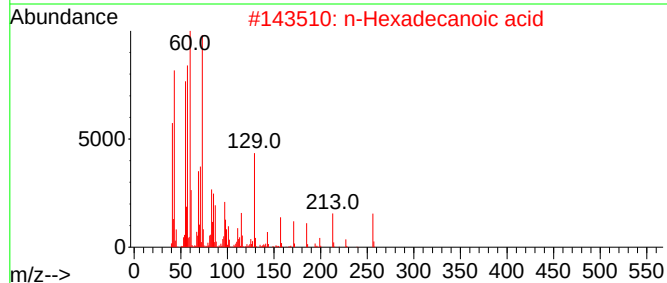
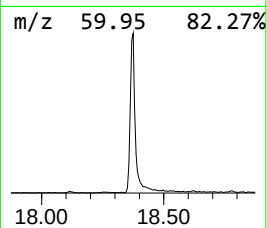
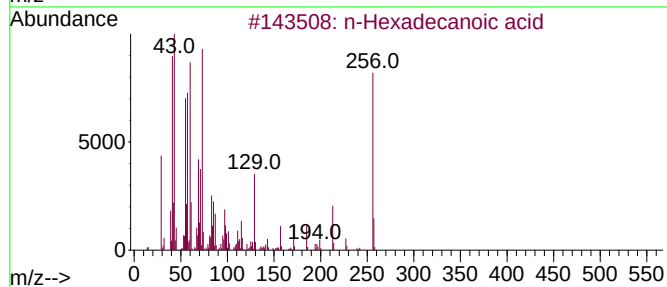
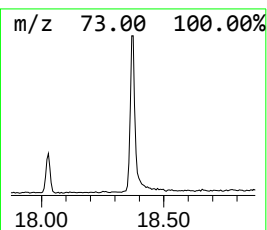
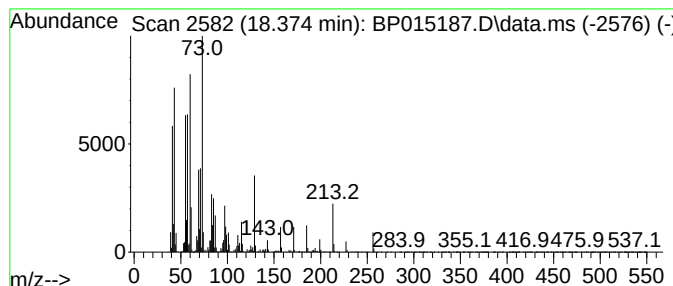
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.374	6.78 ng/ul	1261380	Phenanthrene-d10	17.516

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	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP05187.D
Acq On : 19 May 2023 23:11
Operator : CG/JU
Sample : 02689-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREM3

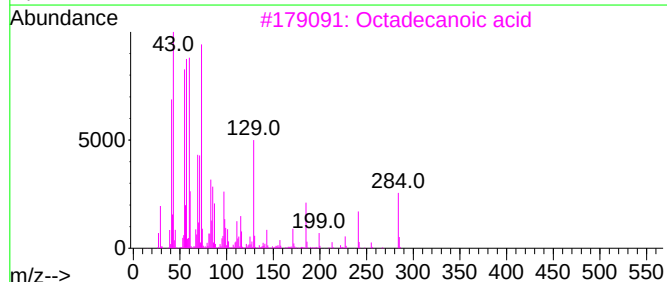
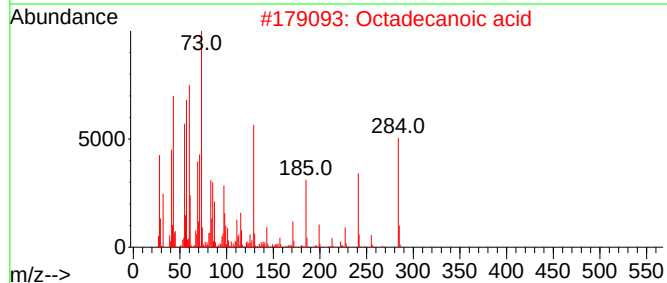
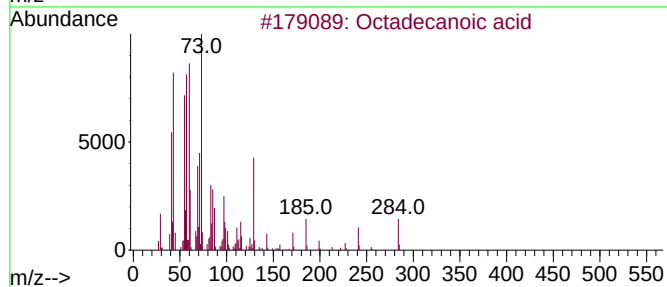
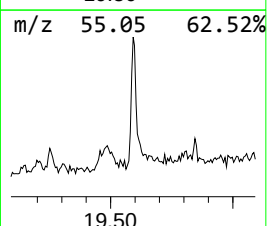
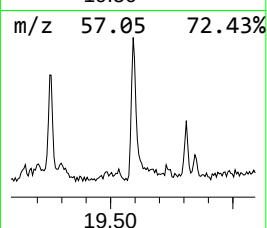
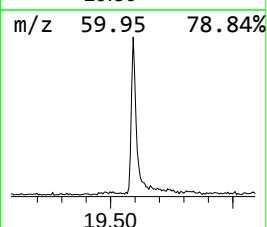
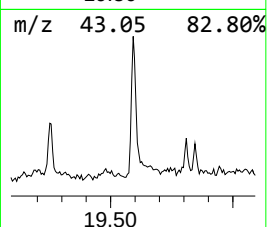
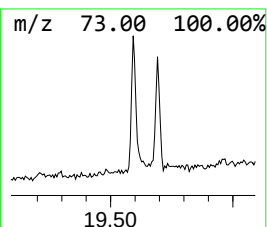
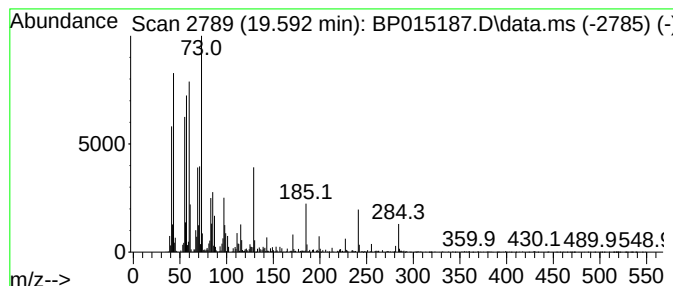
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.592	3.22 ng/ul	468908	Chrysene-d12	21.592

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	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP05187.D
Acq On : 19 May 2023 23:11
Operator : CG/JU
Sample : 02689-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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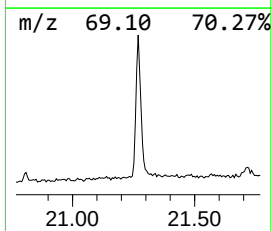
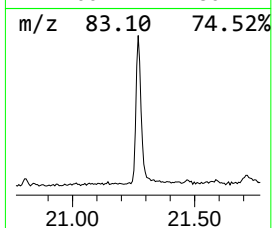
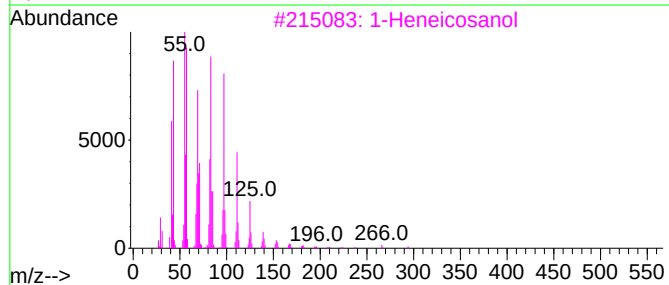
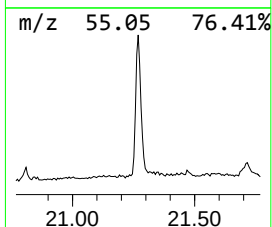
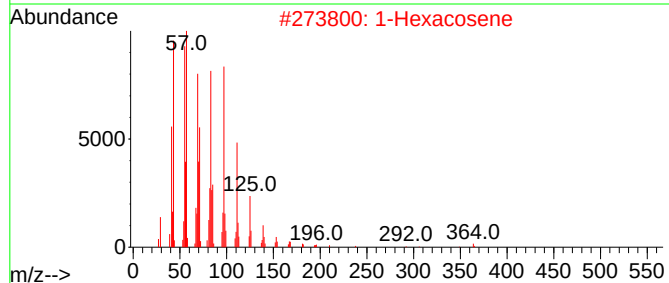
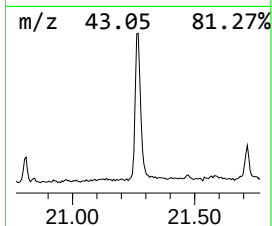
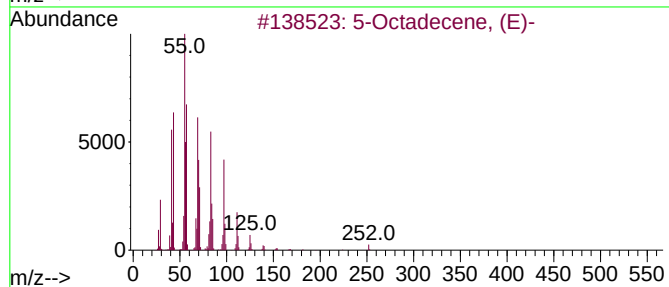
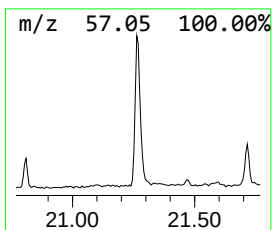
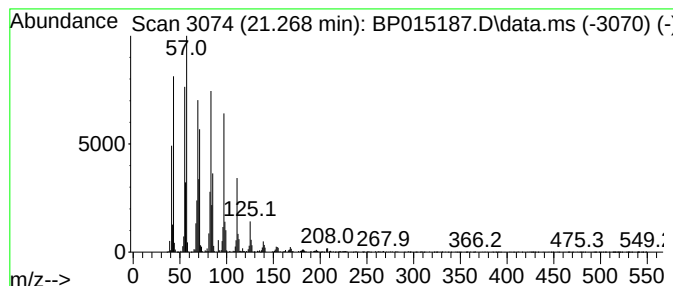
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.268	9.36 ng/ul	1360820	Chrysene-d12	21.592

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2		1-Hexacosene	364	C26H52	018835-33-1	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5		Pentacos-1-ene	350	C25H50	016980-85-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP05187.D
Acq On : 19 May 2023 23:11
Operator : CG/JU
Sample : 02689-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

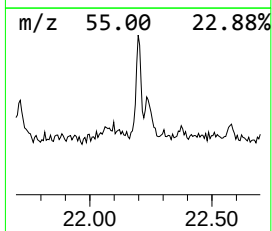
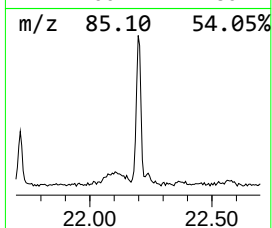
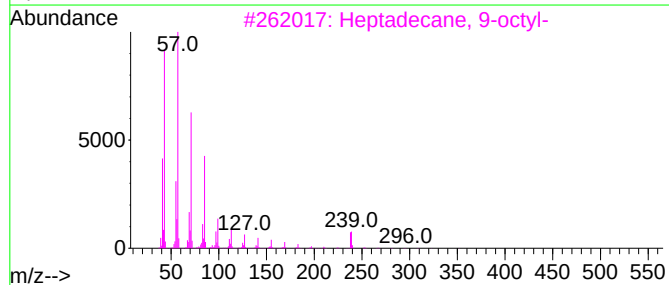
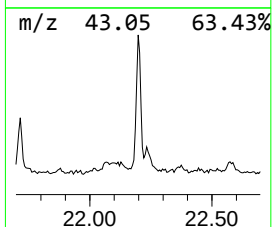
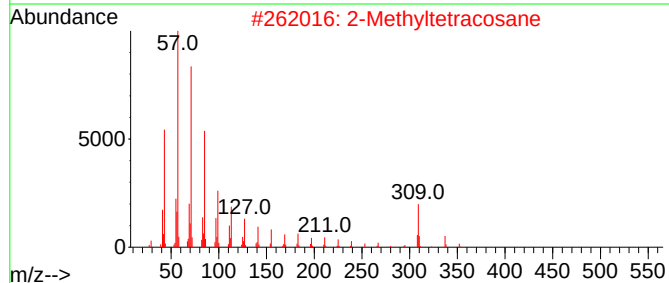
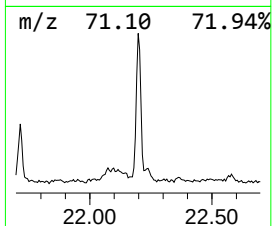
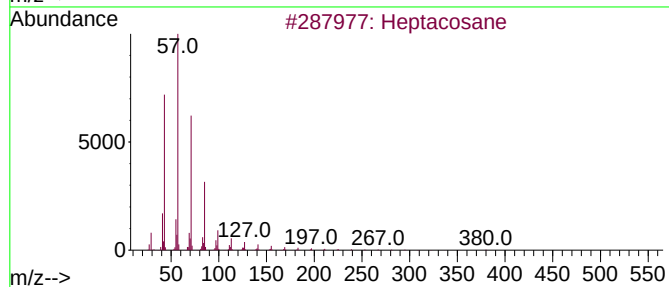
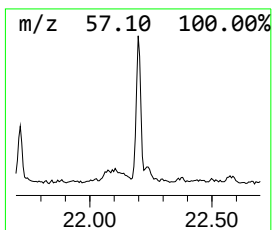
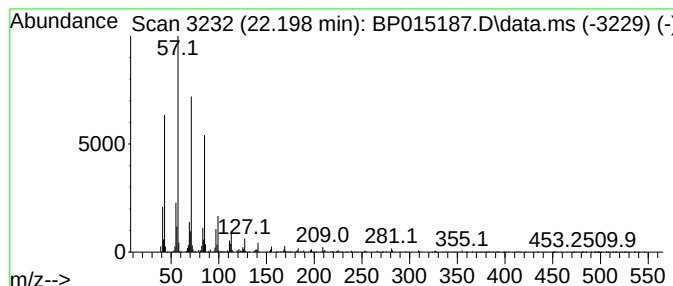
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.198	2.81 ng/ul	409333	Chrysene-d12		21.592	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	95
2	2-Methyltetracosane		352	C25H52	001560-78-7	93
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
4	Tetracosane		338	C24H50	000646-31-1	93
5	Eicosane		282	C20H42	000112-95-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP05187.D
Acq On : 19 May 2023 23:11
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Sample : 02689-11
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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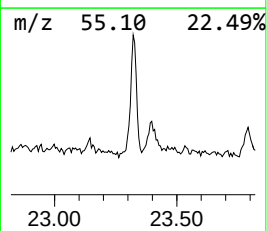
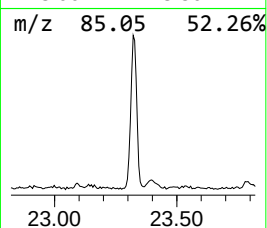
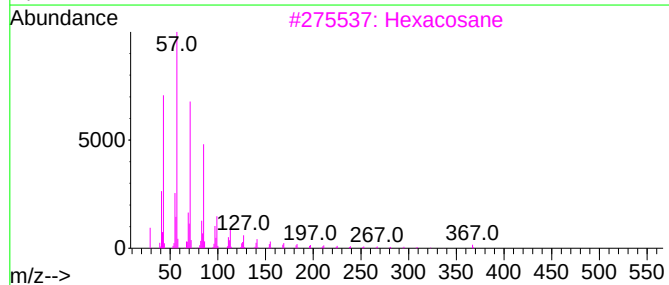
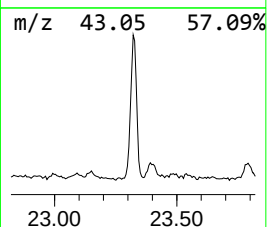
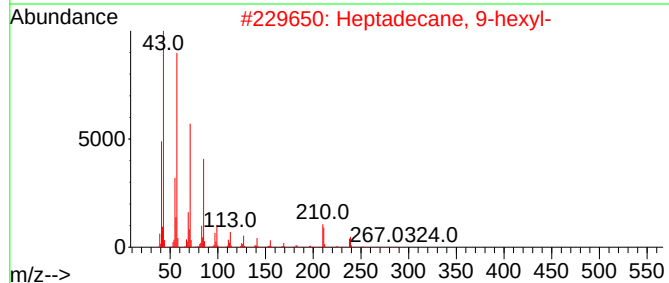
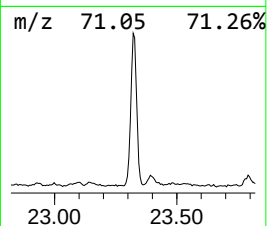
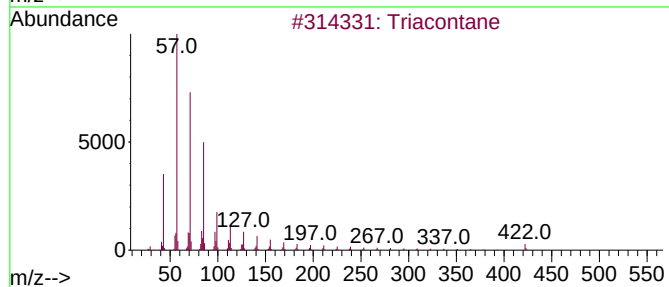
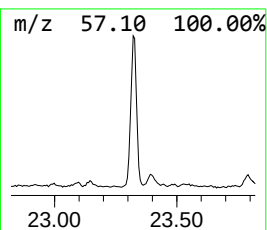
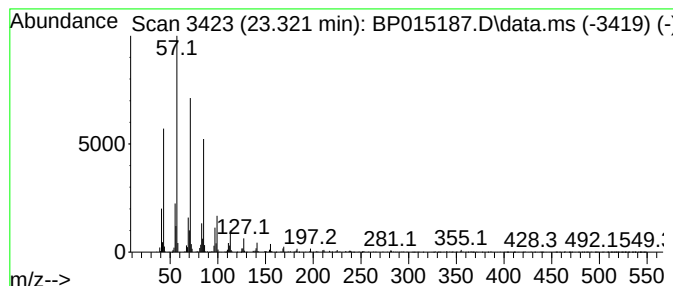
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.321	6.08 ng/ul	832750	Perylene-d12	24.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	96
2			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3			Hexacosane	366	C26H54	000630-01-3	91
4			Eicosane	282	C20H42	000112-95-8	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
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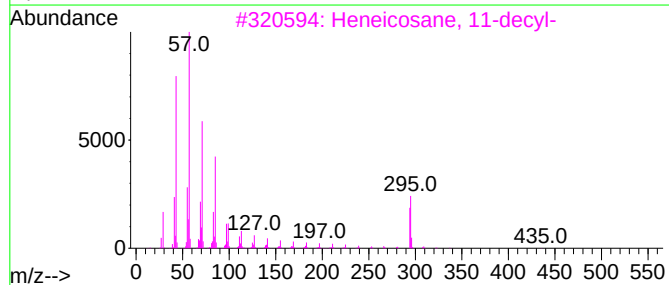
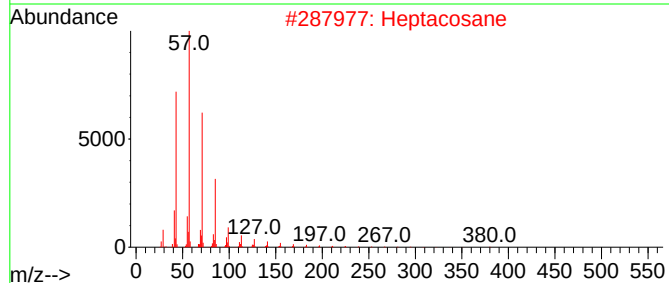
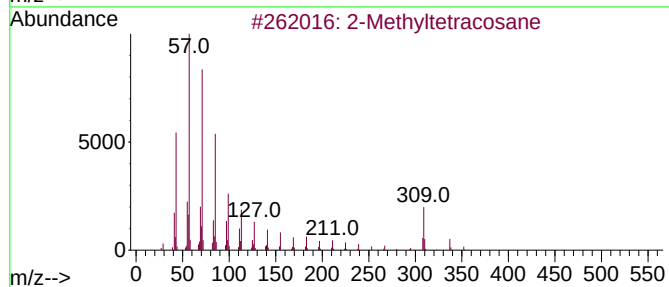
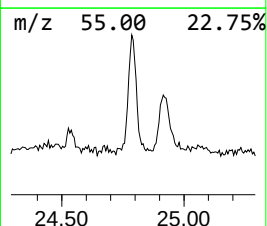
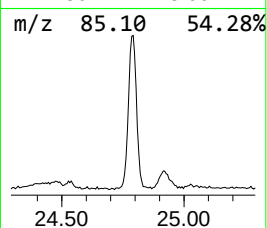
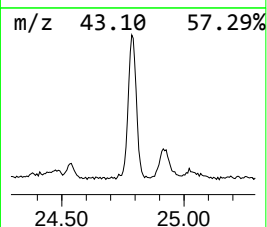
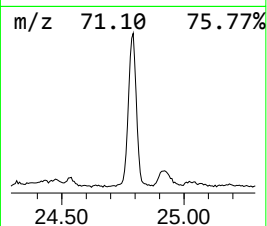
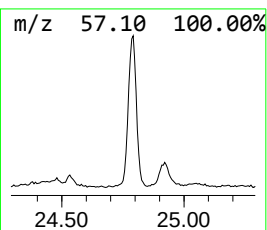
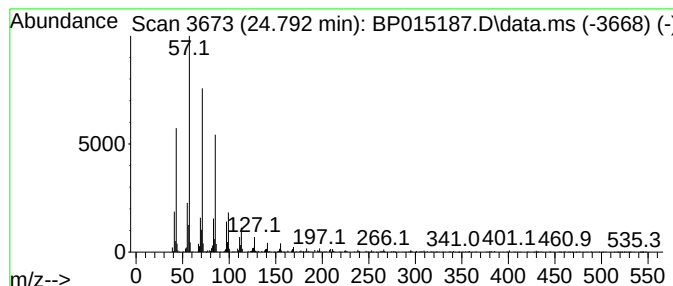
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.792	7.64 ng/ul	1046350	Perylene-d12		24.151	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Methyltetracosane		352	C25H52	001560-78-7	97
2	Heptacosane		380	C27H56	000593-49-7	95
3	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	94
4	Docosane, 5-butyl-		366	C26H54	055282-16-1	94
5	2-Bromo dodecane		248	C12H25Br	013187-99-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP015187.D
Acq On : 19 May 2023 23:11
Operator : CG/JU
Sample : 02689-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREM3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.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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Benzophenone	16.116	4.0	ng/u1	628897	3	14.763	3172770	20.0
n-Hexadecanoic ...	18.374	6.8	ng/u1	1261380	4	17.516	3722470	20.0
Octadecanoic acid	19.592	3.2	ng/u1	468908	5	21.592	2908300	20.0
(DEL) Alkane: S...	21.268	9.4	ng/u1	1360820	5	21.592	2908300	20.0
(DEL) Alkane: S...	22.198	2.8	ng/u1	409333	5	21.592	2908300	20.0
(DEL) Alkane: S...	23.321	6.1	ng/u1	832750	6	24.151	2737520	20.0
(DEL) Alkane: S...	24.792	7.6	ng/u1	1046350	6	24.151	2737520	20.0