

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
 Data File : BN024929.D
 Acq On : 15 Apr 2023 11:37
 Operator : CG/JU
 Sample : 02254-12
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AB4

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_N\Methods\SFAM-EPA-BN041023.M
 Title : SVOA CALIBRATION

Signal : TIC: BN024929.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	9	14	17	rBV	96591	138403	1.41%	0.147%
2	3.305	17	20	22	rVV	72852	98377	1.00%	0.104%
3	3.340	22	26	40	rVB	358038	555291	5.66%	0.589%
4	3.587	65	68	71	rBV4	9417	12549	0.13%	0.013%
5	3.752	94	96	110	rBV	97271	188316	1.92%	0.200%
6	3.970	128	133	140	rBV7	7985	14888	0.15%	0.016%
7	4.070	145	150	157	rBV	22398	36717	0.37%	0.039%
8	4.605	236	241	252	rVB	35514	62969	0.64%	0.067%
9	6.022	478	482	487	rVB6	9489	14598	0.15%	0.015%
10	6.928	629	636	639	rBV8	6474	11758	0.12%	0.012%
11	7.093	657	664	677	rVB	289739	478947	4.88%	0.508%
12	7.264	686	693	707	rVB	1056633	1642069	16.73%	1.743%
13	7.464	719	727	735	rBV	1021817	1626361	16.57%	1.726%
14	7.934	799	807	815	rBV	1211974	1914823	19.51%	2.033%
15	7.999	815	818	822	rVV5	9013	16779	0.17%	0.018%
16	8.052	822	827	833	rVB4	18546	31195	0.32%	0.033%
17	8.158	838	845	849	rBV8	9073	17106	0.17%	0.018%
18	8.646	921	928	938	rBV	824126	1301399	13.26%	1.382%
19	9.105	999	1006	1016	rBV	1311445	2133280	21.73%	2.265%
20	9.210	1020	1024	1029	rVV7	7224	14814	0.15%	0.016%
21	9.263	1029	1033	1040	rVB2	27888	44851	0.46%	0.048%
22	9.510	1069	1075	1080	rBV	14059	23629	0.24%	0.025%
23	9.828	1120	1129	1142	rBV	1080605	1731366	17.64%	1.838%
24	10.058	1165	1168	1173	rBV6	8422	11712	0.12%	0.012%
25	10.363	1213	1220	1235	rBV	1401305	2349244	23.94%	2.494%
26	10.528	1243	1248	1256	rBV3	31406	64633	0.66%	0.069%
27	10.746	1277	1285	1297	rVB	1696641	2820770	28.74%	2.994%
28	10.840	1297	1301	1303	rBV4	9867	15648	0.16%	0.017%
29	10.887	1303	1309	1322	rVB	857565	1472417	15.00%	1.563%
30	11.281	1373	1376	1383	rVB6	8048	13151	0.13%	0.014%
31	11.352	1383	1388	1391	rBV4	7452	12191	0.12%	0.013%
32	11.534	1414	1419	1424	rBV2	16548	30290	0.31%	0.032%
33	12.010	1495	1500	1507	rVB9	6793	13733	0.14%	0.015%
34	12.334	1551	1555	1560	rVB	24054	38506	0.39%	0.041%
35	12.575	1588	1596	1599	rBV9	8159	14085	0.14%	0.015%
36	13.399	1732	1736	1740	rVB3	11573	16277	0.17%	0.017%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
 Title : SVOA CALIBRATION

37	13.540	1756	1760	1767	rVB9	8032	14377	0.15%	0.015%
38	13.781	1796	1801	1808	rBV3	34025	52051	0.53%	0.055%
39	13.987	1829	1836	1847	rBV	2798220	3936731	40.11%	4.179%
40	14.104	1851	1856	1860	rVB7	10567	15374	0.16%	0.016%
41	14.275	1878	1885	1896	rBV	3171993	4537767	46.23%	4.817%
42	14.469	1914	1918	1922	rVB2	21513	26932	0.27%	0.029%
43	14.581	1928	1937	1948	rBV	2605831	3747122	38.18%	3.978%
44	14.775	1962	1970	1988	rBV	215988	425884	4.34%	0.452%
45	14.987	2000	2006	2010	rBV6	16444	25925	0.26%	0.028%
46	15.081	2019	2022	2027	rVB3	15226	20925	0.21%	0.022%
47	15.157	2030	2035	2038	rBV4	8955	12033	0.12%	0.013%
48	15.316	2057	2062	2066	rBV5	11534	17444	0.18%	0.019%
49	15.416	2072	2079	2084	rVV	186657	244613	2.49%	0.260%
50	15.569	2099	2105	2115	rBV	4171316	5920406	60.32%	6.285%
51	15.687	2119	2125	2138	rVV	1674889	2278824	23.22%	2.419%
52	15.822	2139	2148	2153	rVV3	93583	186076	1.90%	0.198%
53	15.863	2153	2155	2158	rVV2	14650	15303	0.16%	0.016%
54	15.934	2161	2167	2172	rVV	629470	894703	9.12%	0.950%
55	15.975	2172	2174	2180	rVV3	42257	61929	0.63%	0.066%
56	16.040	2181	2185	2190	rVB2	44201	58920	0.60%	0.063%
57	16.281	2221	2226	2234	rBV	638047	1049717	10.69%	1.114%
58	16.498	2260	2263	2267	rBV5	16355	22667	0.23%	0.024%
59	16.587	2276	2278	2280	rBV2	10450	10812	0.11%	0.011%
60	16.622	2280	2284	2288	rVV2	55216	95695	0.97%	0.102%
61	16.687	2292	2295	2298	rVV	41703	49016	0.50%	0.052%
62	16.740	2301	2304	2308	rVB	39530	44850	0.46%	0.048%
63	16.787	2308	2312	2318	rBV	60138	87525	0.89%	0.093%
64	16.857	2321	2324	2327	rVV	41457	43372	0.44%	0.046%
65	17.075	2356	2361	2365	rBV	803173	955224	9.73%	1.014%
66	17.128	2365	2370	2374	rVB	225531	312304	3.18%	0.332%
67	17.328	2398	2404	2410	rBV	3273217	4614934	47.02%	4.899%
68	17.428	2415	2421	2429	rVB	5010142	7161970	72.97%	7.603%
69	17.551	2438	2442	2445	rBV	64699	86008	0.88%	0.091%
70	17.616	2448	2453	2459	rBV3	88642	165808	1.69%	0.176%
71	17.681	2461	2464	2469	rBV5	37831	64895	0.66%	0.069%
72	17.734	2470	2473	2481	rVV2	67968	116049	1.18%	0.123%
73	17.816	2482	2487	2493	rVV	827425	973934	9.92%	1.034%
74	17.992	2513	2517	2522	rVB	404257	481246	4.90%	0.511%
75	18.098	2532	2535	2540	rBV2	49402	83819	0.85%	0.089%
76	18.216	2550	2555	2560	rBV	596227	821218	8.37%	0.872%

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 Title : SVOA CALIBRATION

77	18.510	2601	2605	2610	rBV	703010	806907	8.22%	0.857%
78	19.139	2708	2712	2715	rBV	529358	682297	6.95%	0.724%
79	19.298	2735	2739	2744	rVB2	191990	257090	2.62%	0.273%
80	19.357	2746	2749	2754	rVB3	87264	128482	1.31%	0.136%
81	19.492	2768	2772	2781	rBV	278641	418092	4.26%	0.444%
82	19.722	2805	2811	2824	rVB	6816695	8966391	91.35%	9.518%
83	20.257	2898	2902	2905	rVB	221936	262800	2.68%	0.279%
84	20.751	2982	2986	2990	rVB	162272	192454	1.96%	0.204%
85	21.222	3061	3066	3075	rBV	665607	1187675	12.10%	1.261%
86	21.516	3111	3116	3130	rVB	4180927	5377313	54.79%	5.708%
87	22.422	3266	3270	3277	rBV	570763	759431	7.74%	0.806%
88	23.204	3399	3403	3410	rVB2	215287	325889	3.32%	0.346%
89	23.816	3499	3507	3520	rBV2	4672250	9815033	100.00%	10.419%
90	23.969	3526	3533	3551	rVB2	2761267	5750207	58.59%	6.104%
91	24.586	3634	3638	3647	rVB3	274262	553474	5.64%	0.588%

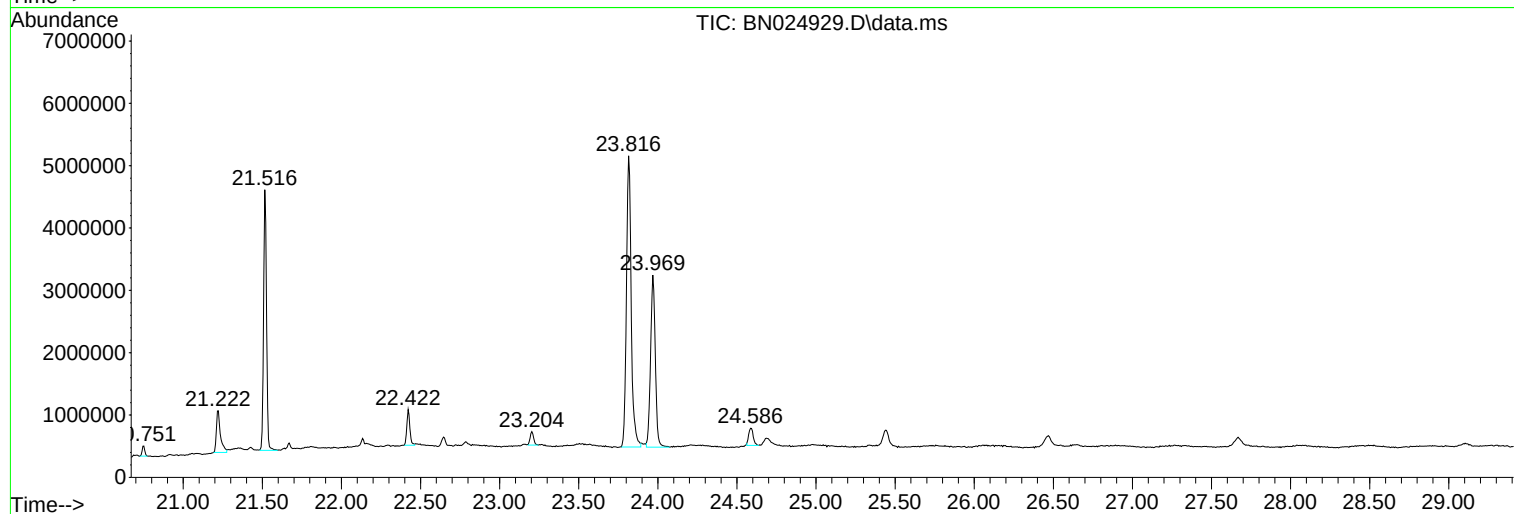
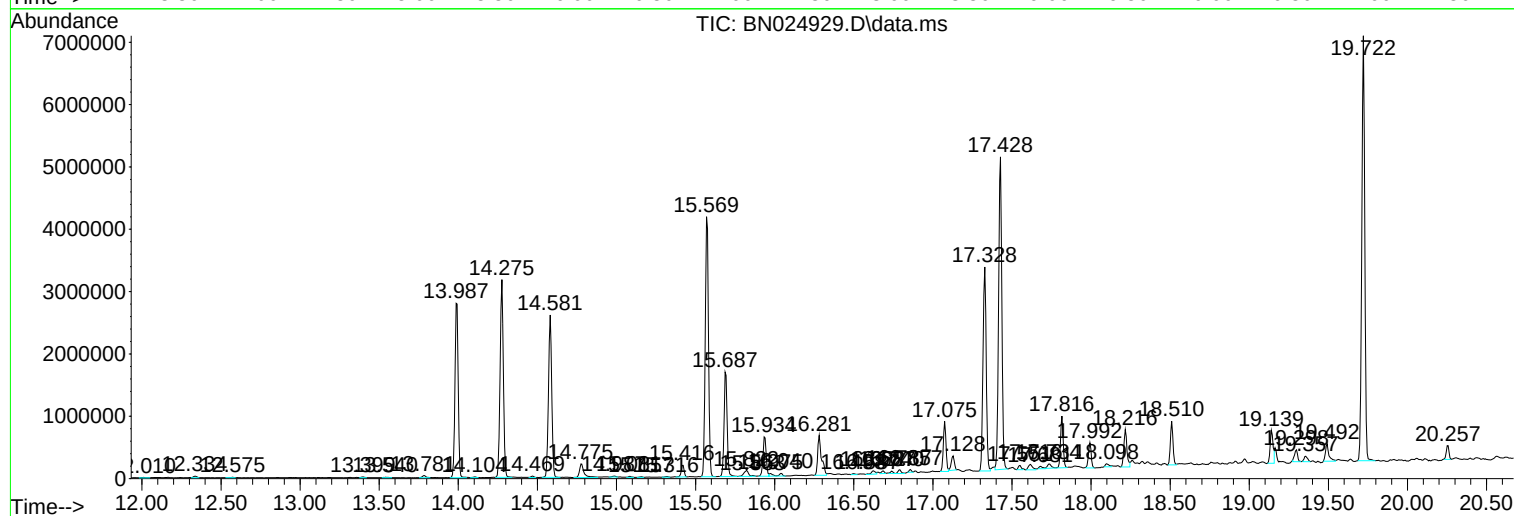
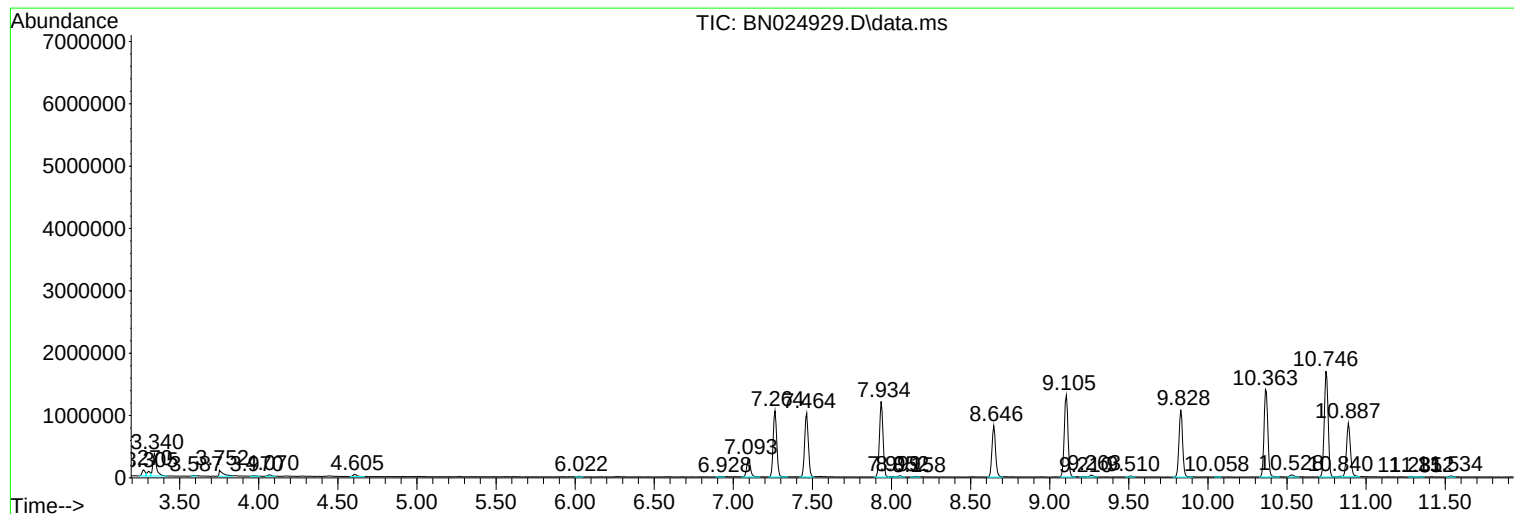
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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



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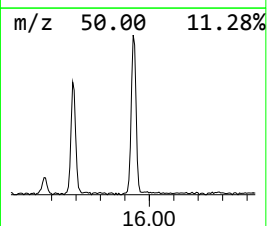
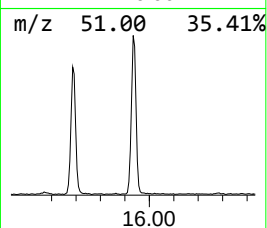
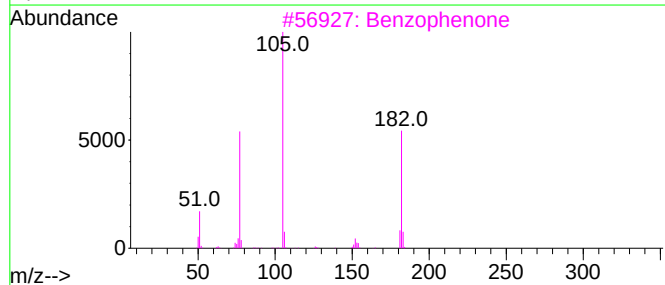
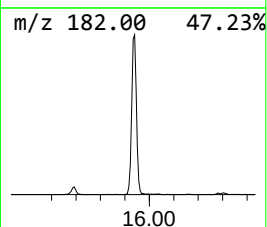
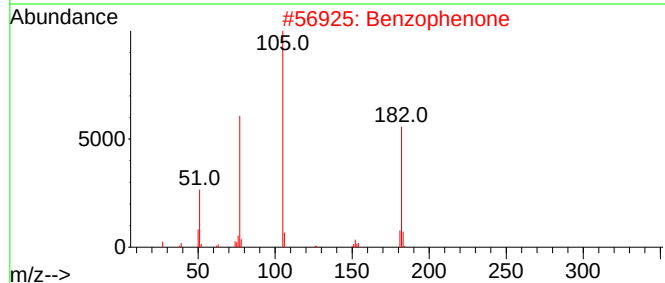
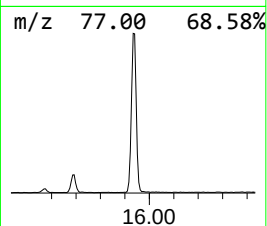
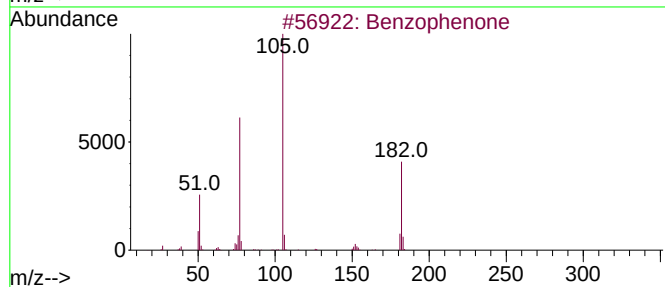
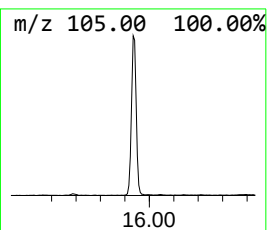
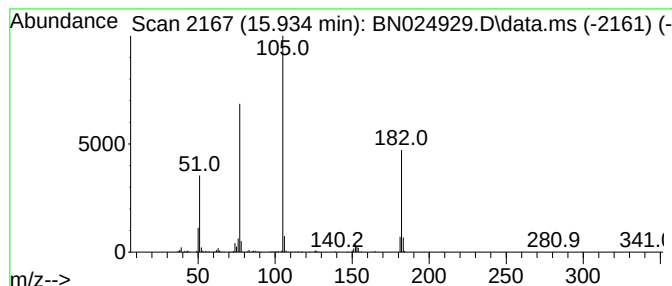
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.934	4.78 ng/ul	894703	Acenaphthene-d10	14.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	97
3			Benzophenone	182	C13H10O	000119-61-9	97
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



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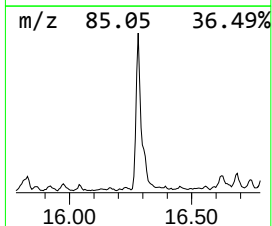
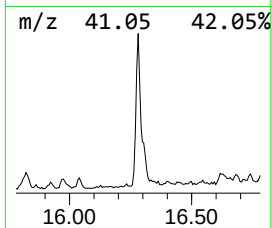
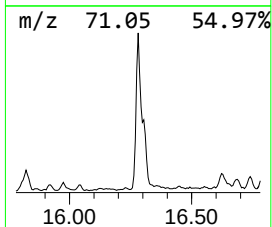
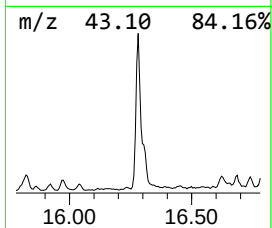
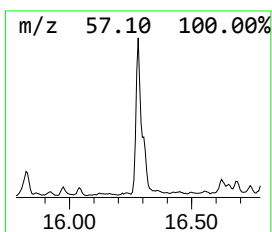
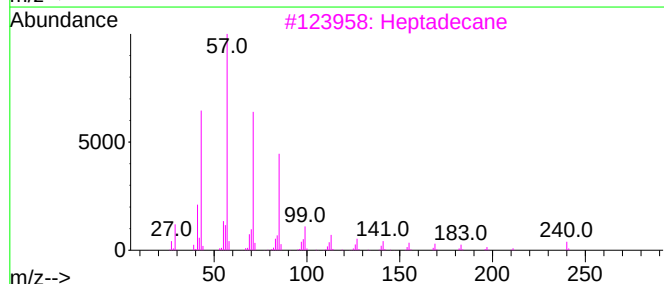
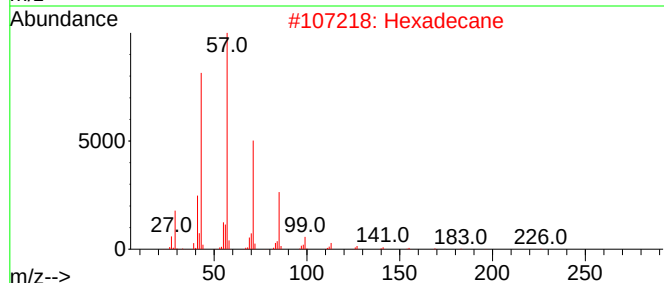
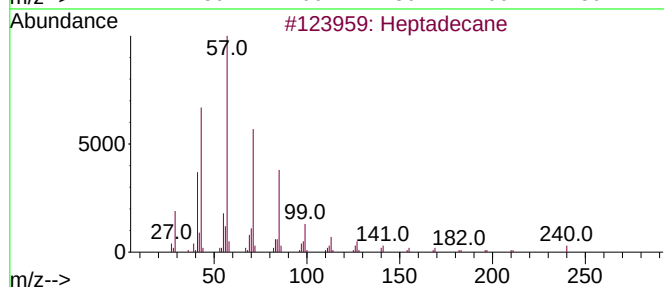
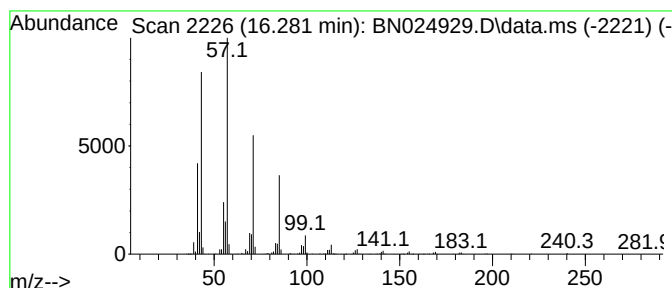
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.281	4.55 ng/ul	1049720	Phenanthrene-d10	17.328	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Hexadecane	226	C16H34	000544-76-3	91
3	Heptadecane	240	C17H36	000629-78-7	90
4	Heneicosane	296	C21H44	000629-94-7	90
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90



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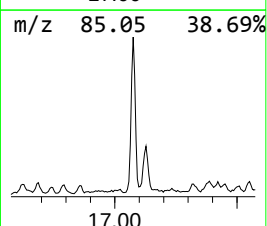
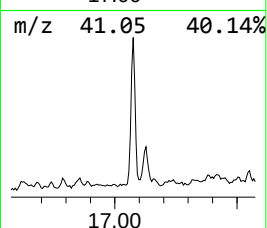
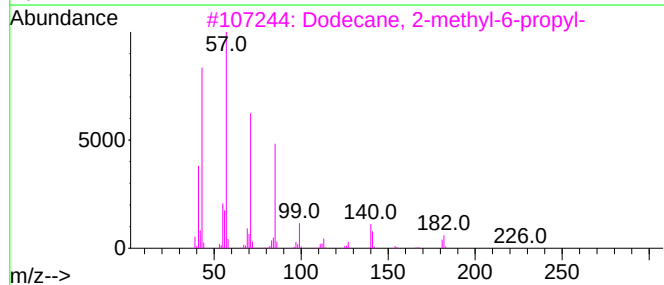
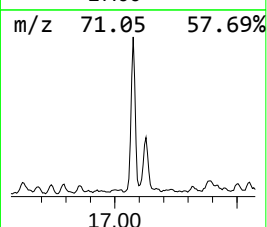
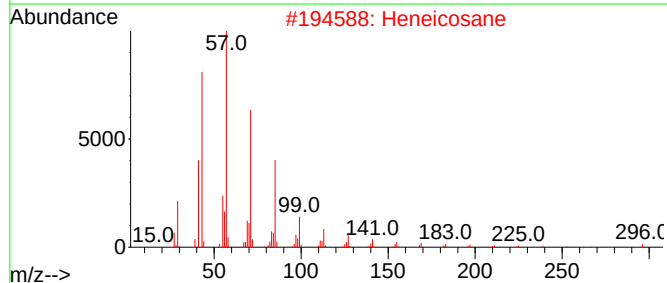
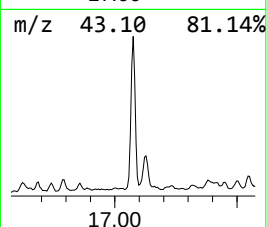
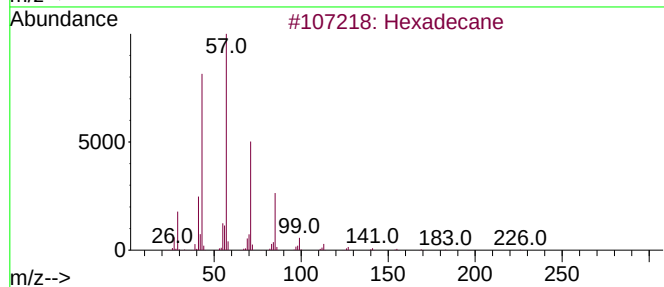
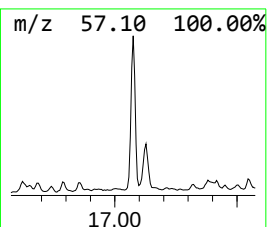
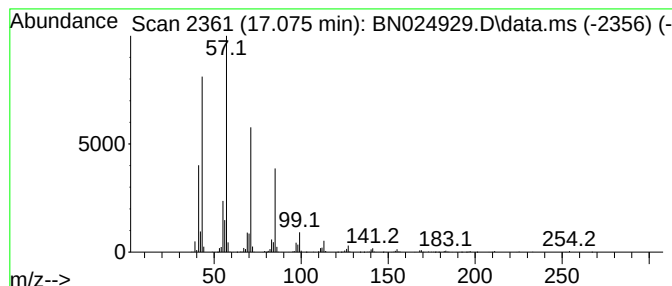
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.075	4.14 ng/ul	955224	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4			Eicosane	282	C20H42	000112-95-8	90
5			Heptadecane	240	C17H36	000629-78-7	86



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ClientSampleId :
C0AB4

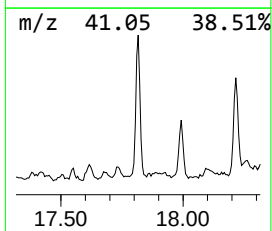
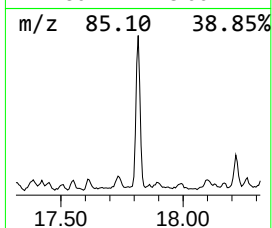
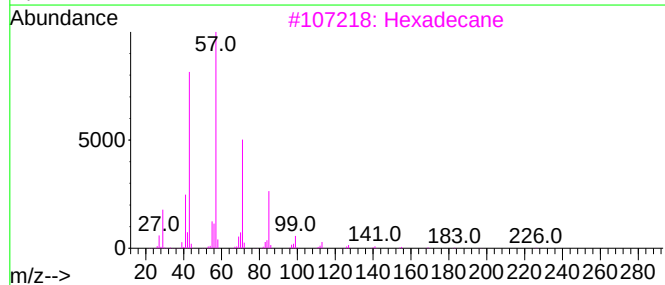
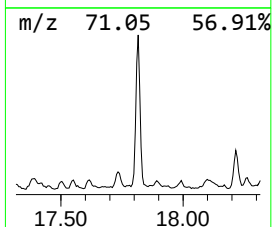
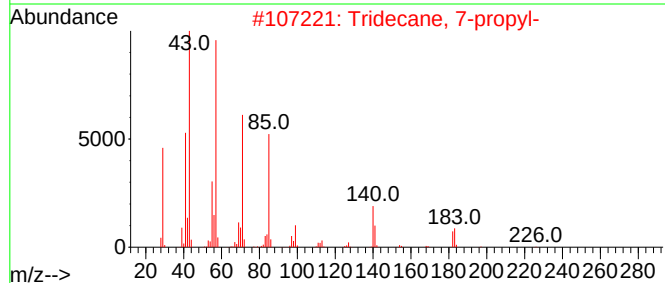
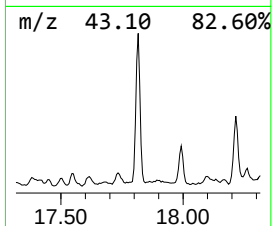
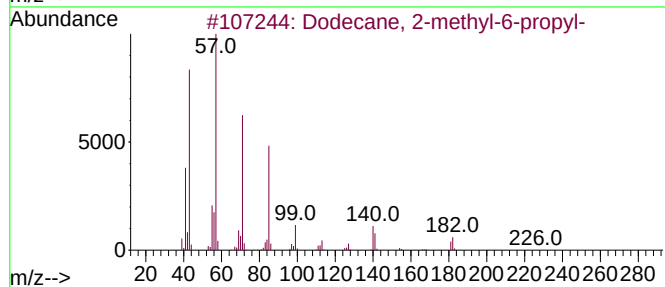
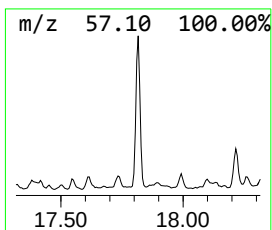
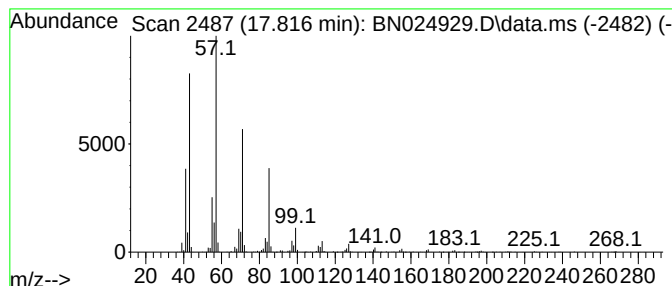
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.816	4.22 ng/ul	973934	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
2			Tridecane, 7-propyl-	226	C16H34	055045-09-5	91
3			Hexadecane	226	C16H34	000544-76-3	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024929.D
Acq On : 15 Apr 2023 11:37
Operator : CG/JU
Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
BNA_N
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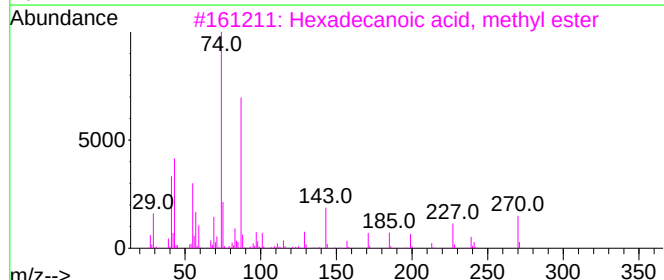
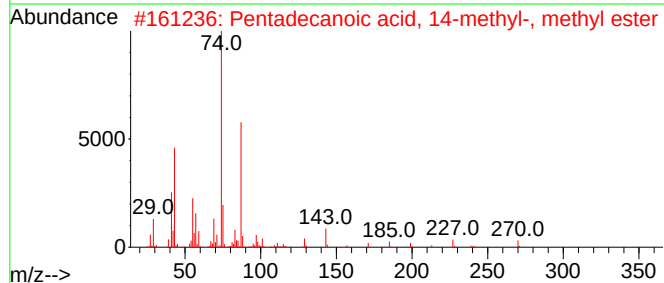
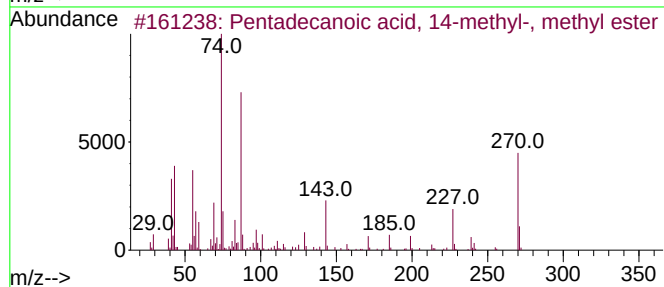
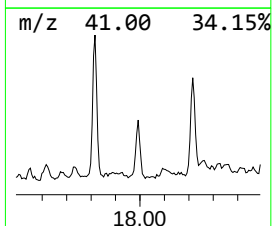
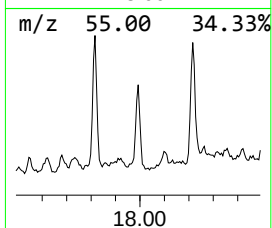
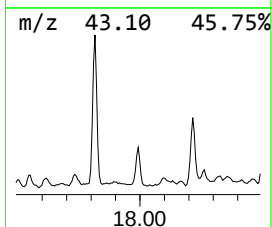
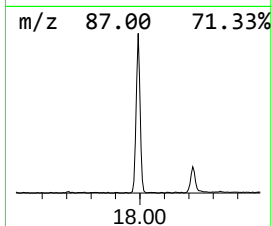
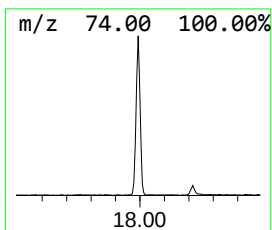
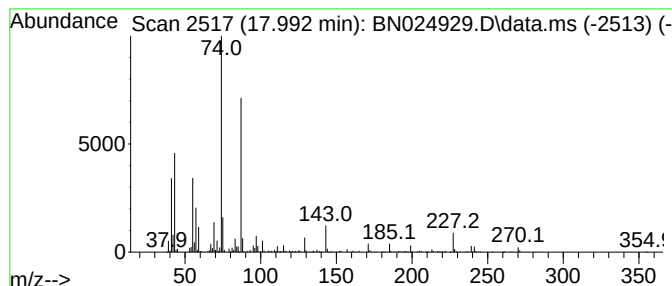
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecanoic acid, 14-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	2.09 ng/ul	481246	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024929.D
Acq On : 15 Apr 2023 11:37
Operator : CG/JU
Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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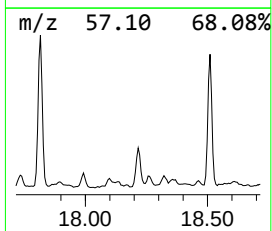
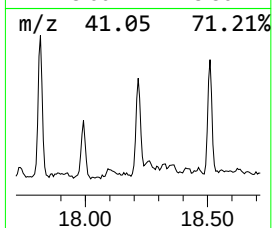
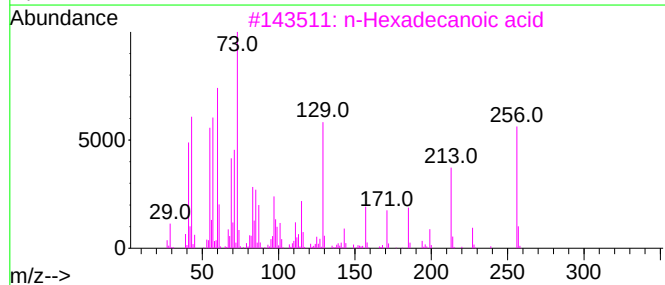
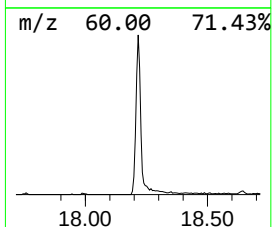
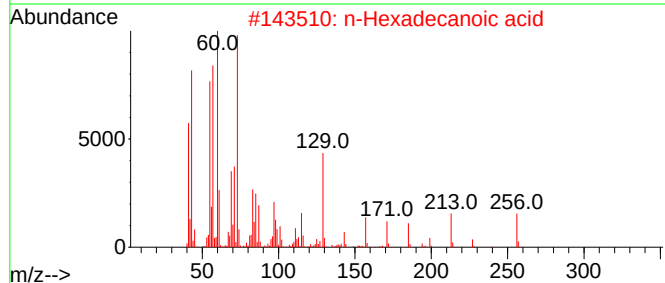
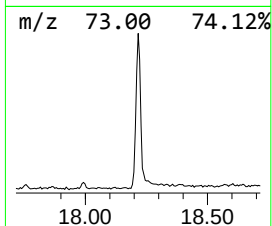
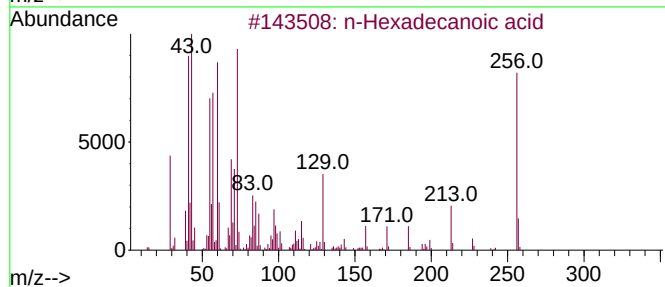
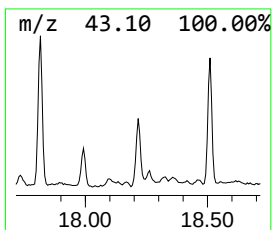
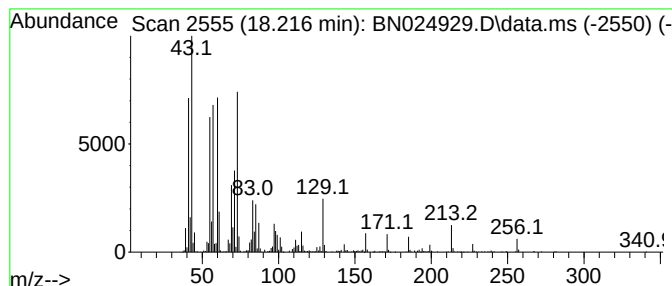
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	3.56 ng/ul	821218	Phenanthrene-d10	17.328

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	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	91
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024929.D
Acq On : 15 Apr 2023 11:37
Operator : CG/JU
Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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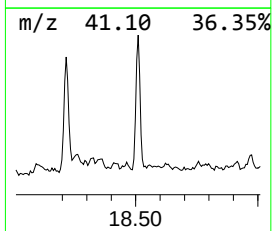
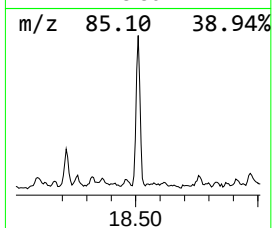
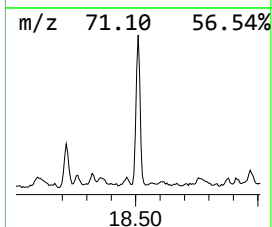
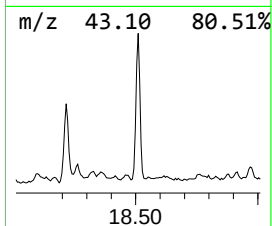
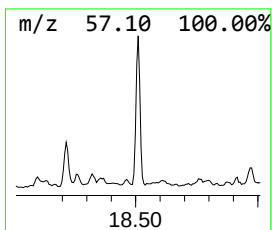
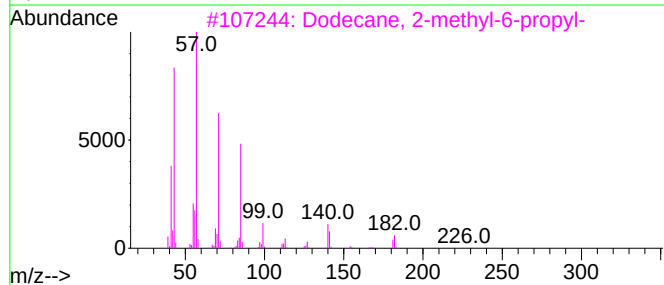
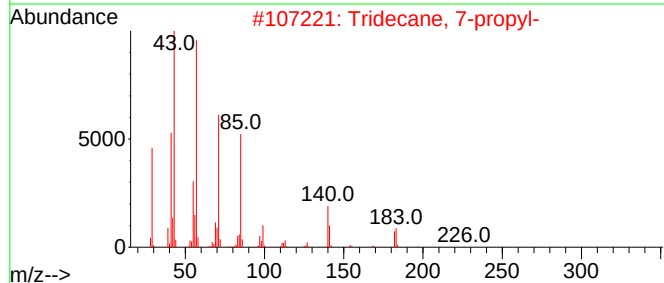
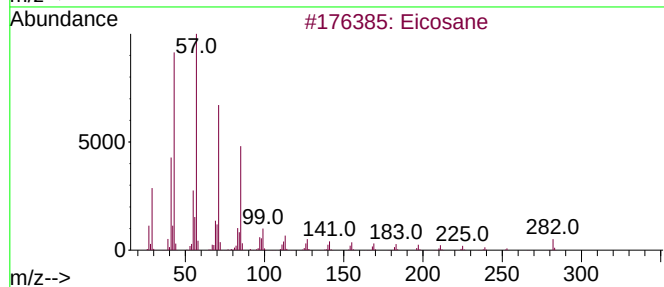
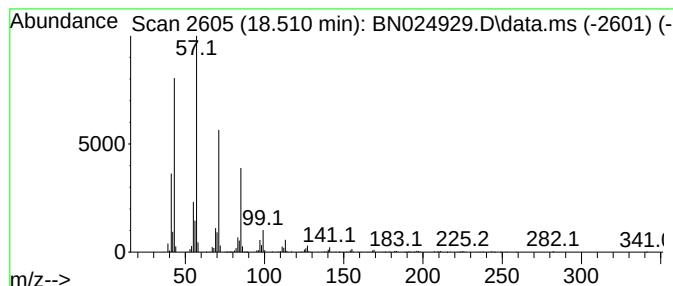
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.510	3.50 ng/ul	806907	Phenanthrene-d10	17.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Tridecane, 7-propyl-	226	C16H34	055045-09-5	93
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
4			Nonadecane	268	C19H40	000629-92-5	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024929.D
Acq On : 15 Apr 2023 11:37
Operator : CG/JU
Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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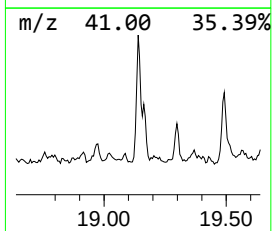
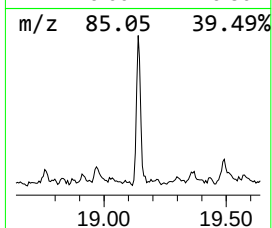
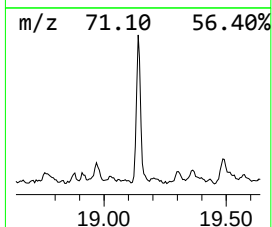
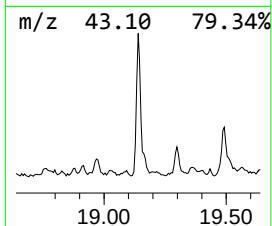
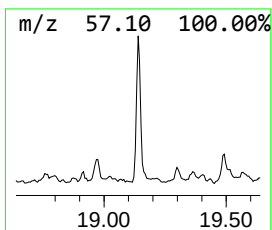
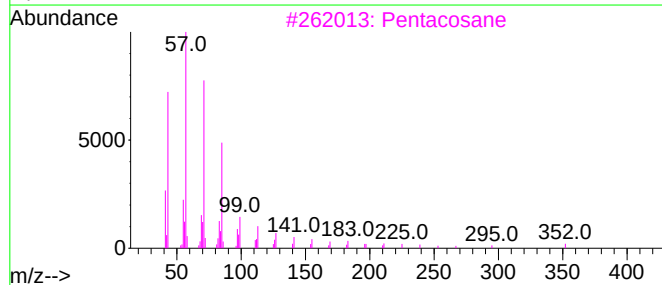
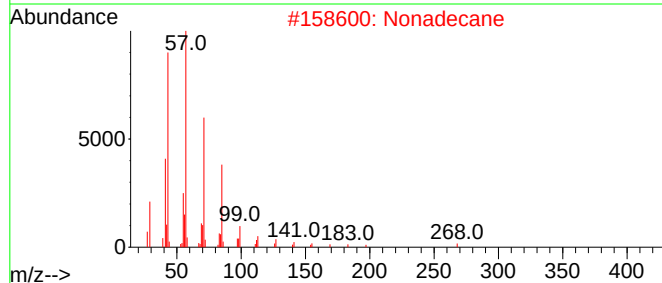
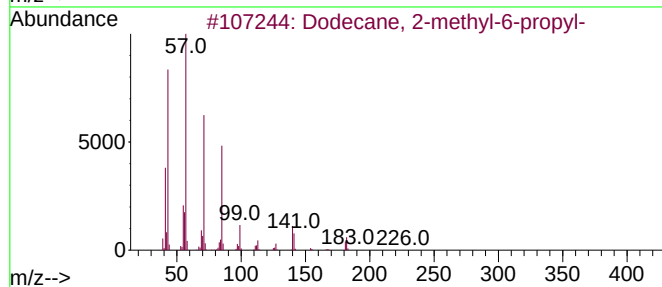
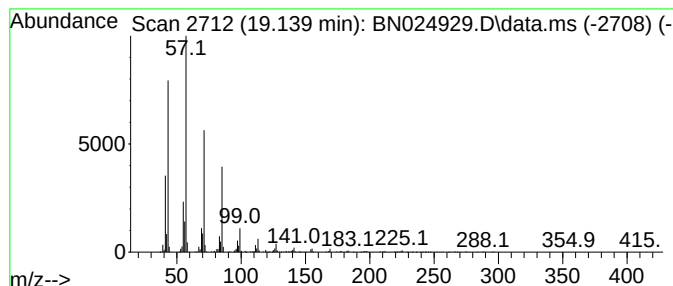
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	2.96 ng/ul	682297	Phenanthrene-d10	17.328

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	94
2		Nonadecane	268	C19H40	000629-92-5	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
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Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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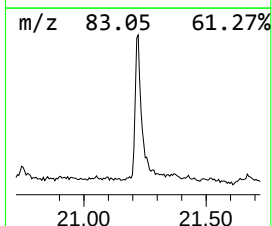
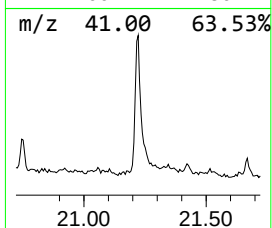
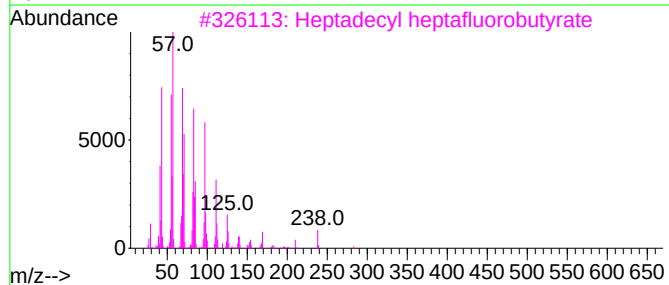
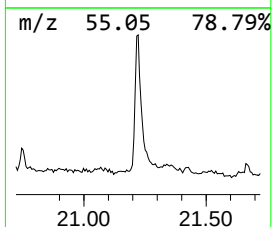
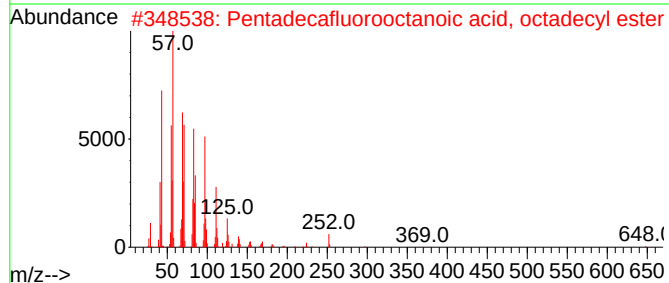
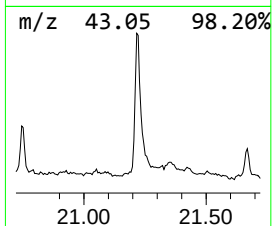
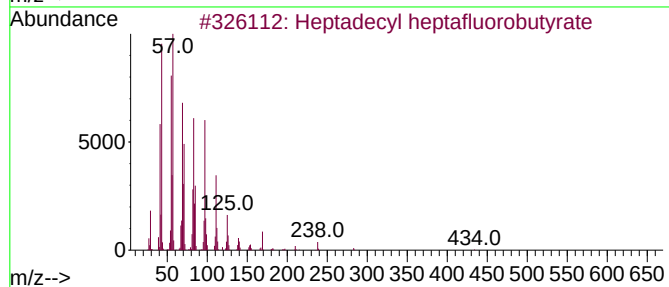
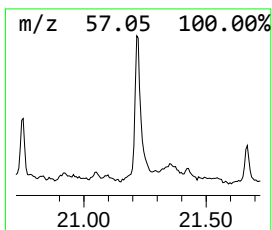
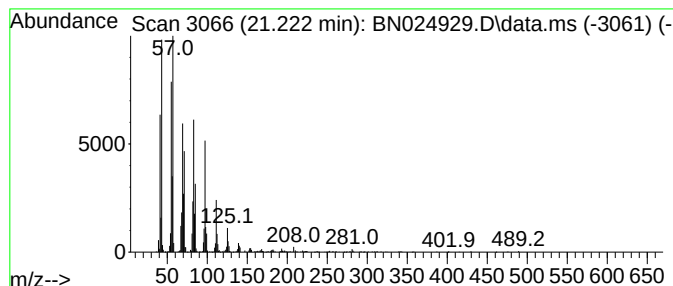
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	4.42 ng/ul	1187680	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
5			1-Docosene	308	C22H44	001599-67-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
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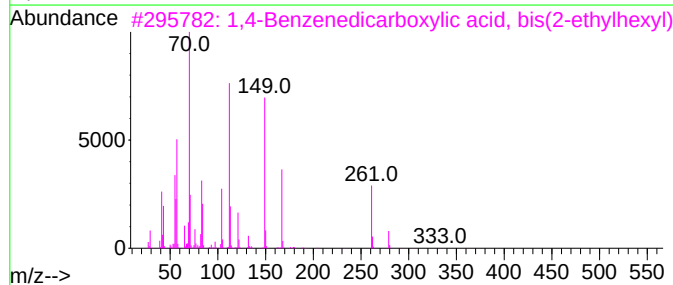
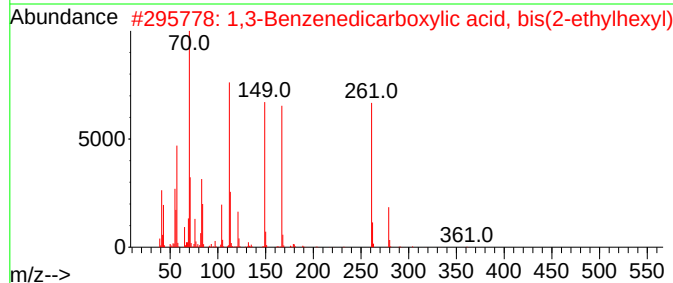
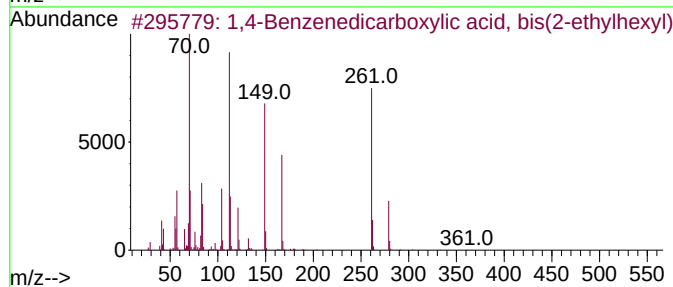
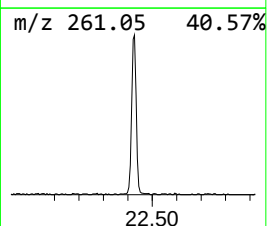
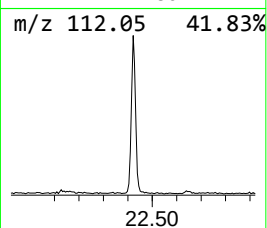
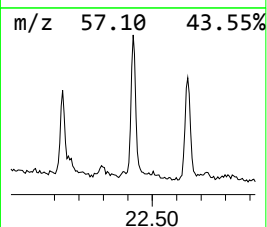
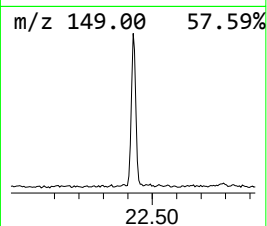
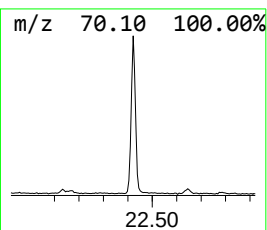
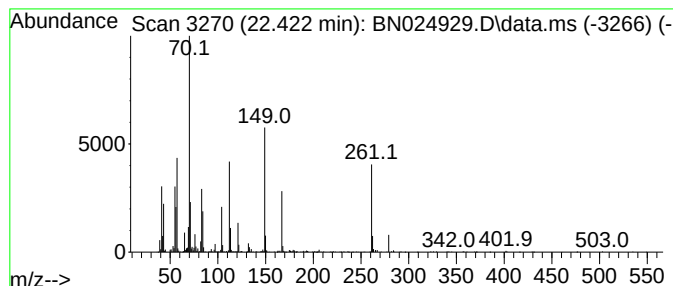
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,4-Benzenedicarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.422	2.82 ng/ul	759431	Chrysene-d12	21.516	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	59
2	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	58
3	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	58
4	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	53
5	1-(2-Hydroxymethylpyrrolidin-1-y...	143	C7H13NO2	1000192-83-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041423\
Data File : BN024929.D
Acq On : 15 Apr 2023 11:37
Operator : CG/JU
Sample : 02254-12
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AB4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.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
Benzophenone	15.934	4.8	ng/ul	894703	3	14.581	3747120	20.0
(DEL) Alkane: S...	16.281	4.5	ng/ul	1049720	4	17.328	4614930	20.0
(DEL) Alkane: S...	17.075	4.1	ng/ul	955224	4	17.328	4614930	20.0
(DEL) Alkane: S...	17.816	4.2	ng/ul	973934	4	17.328	4614930	20.0
Pentadecanoic a...	17.992	2.1	ng/ul	481246	4	17.328	4614930	20.0
n-Hexadecanoic ...	18.216	3.6	ng/ul	821218	4	17.328	4614930	20.0
(DEL) Alkane: S...	18.510	3.5	ng/ul	806907	4	17.328	4614930	20.0
(DEL) Alkane: S...	19.139	3.0	ng/ul	682297	4	17.328	4614930	20.0
Heptadecyl hept...	21.222	4.4	ng/ul	1187680	5	21.516	5377310	20.0
1,4-Benzenedica...	22.422	2.8	ng/ul	759431	5	21.516	5377310	20.0