

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015079.D
 Acq On : 25 Jun 2021 15:55
 Operator : CG/JU
 Sample : M2680-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD34

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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-BN061821MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015079.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	31	34	40	rVB	43855	57950	1.38%	0.128%
2	3.681	97	101	113	rBV	331351	522646	12.48%	1.153%
3	3.787	116	119	125	rVB3	23596	34944	0.83%	0.077%
4	3.911	138	140	146	rVB3	6171	7428	0.18%	0.016%
5	4.005	151	156	165	rVB	43736	69932	1.67%	0.154%
6	4.358	212	216	221	rVB3	16480	24485	0.58%	0.054%
7	4.487	234	238	242	rBV2	12155	15189	0.36%	0.034%
8	4.534	242	246	252	rVB	44926	63157	1.51%	0.139%
9	5.093	337	341	347	rVB	14960	21235	0.51%	0.047%
10	5.287	369	374	378	rBV	34467	47667	1.14%	0.105%
11	5.411	390	395	404	rVB	96488	195630	4.67%	0.432%
12	5.775	451	457	464	rVB	47410	73201	1.75%	0.161%
13	6.746	618	622	627	rVB4	5382	9509	0.23%	0.021%
14	6.893	640	647	656	rBV	645090	1020253	24.37%	2.251%
15	7.063	668	676	686	rBV	512278	814434	19.45%	1.797%
16	7.263	703	710	720	rVV	666933	1067097	25.49%	2.354%
17	7.363	723	727	734	rVB8	6606	12701	0.30%	0.028%
18	7.728	782	789	799	rBV	565952	910021	21.73%	2.008%
19	7.963	824	829	837	rVB2	8480	15632	0.37%	0.034%
20	8.075	842	848	852	rBV	13980	21666	0.52%	0.048%
21	8.275	876	882	889	rBV4	6941	13772	0.33%	0.030%
22	8.416	899	906	921	rBV	750807	1225410	29.27%	2.703%
23	8.540	921	927	933	rVB2	18535	32271	0.77%	0.071%
24	8.816	969	974	977	rBV2	10814	16954	0.40%	0.037%
25	8.869	977	983	990	rVV	579476	938693	22.42%	2.071%
26	8.963	994	999	1005	rVB6	2912	7634	0.18%	0.017%
27	9.322	1054	1060	1066	rBV2	3938	8075	0.19%	0.018%
28	9.587	1094	1105	1113	rBV	415126	672823	16.07%	1.484%
29	10.116	1188	1195	1204	rBV	741032	1202746	28.73%	2.653%
30	10.275	1215	1222	1237	rBV	217417	364279	8.70%	0.804%
31	10.498	1253	1260	1268	rBV	753371	1217435	29.08%	2.686%
32	10.628	1274	1282	1293	rBV	693361	1192929	28.49%	2.632%
33	12.087	1525	1530	1536	rVB2	13583	19223	0.46%	0.042%
34	12.710	1631	1636	1641	rVB4	5964	9090	0.22%	0.020%
35	12.963	1673	1679	1683	rVB	9307	13688	0.33%	0.030%
36	13.187	1713	1717	1724	rBV	14247	18658	0.45%	0.041%

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Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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-BN061821MA.M
 Title : SVOA CALIBRATION

37	13.763	1809	1815	1821	rBV	1270796	1706917	40.77%	3.765%
38	14.039	1854	1862	1871	rBV	1427775	2182042	52.11%	4.814%
39	14.269	1896	1901	1905	rBV	19344	26028	0.62%	0.057%
40	14.351	1909	1915	1925	rBV	1047092	1539784	36.78%	3.397%

41	14.475	1929	1936	1939	rBV2	5031	8100	0.19%	0.018%
42	14.539	1939	1947	1959	rBV	512033	716914	17.12%	1.582%
43	14.769	1982	1986	1991	rVB	27945	34692	0.83%	0.077%
44	15.175	2049	2055	2058	rBV5	5356	9905	0.24%	0.022%
45	15.222	2058	2063	2068	rVB	14527	19139	0.46%	0.042%

46	15.345	2078	2084	2091	rVB	1890977	2661083	63.56%	5.870%
47	15.457	2098	2103	2109	rBV	389293	526628	12.58%	1.162%
48	15.510	2109	2112	2118	rVB	29305	39376	0.94%	0.087%
49	15.586	2121	2125	2129	rBV2	20922	28286	0.68%	0.062%
50	15.775	2151	2157	2164	rBV	183241	234910	5.61%	0.518%

51	16.086	2207	2210	2212	rBV2	10509	12869	0.31%	0.028%
52	16.139	2216	2219	2224	rVB5	14444	23262	0.56%	0.051%
53	16.251	2234	2238	2244	rBV6	6261	10300	0.25%	0.023%
54	16.345	2250	2254	2260	rBV8	4647	8659	0.21%	0.019%
55	16.851	2333	2340	2342	rBV4	16647	28977	0.69%	0.064%

56	16.881	2342	2345	2348	rVV2	26688	31902	0.76%	0.070%
57	16.933	2349	2354	2357	rVV3	7424	12634	0.30%	0.028%
58	17.098	2376	2382	2387	rBV	1254536	1786886	42.68%	3.942%
59	17.139	2387	2389	2393	rVV	43254	53262	1.27%	0.117%
60	17.198	2393	2399	2410	rVV	1986456	2943283	70.30%	6.493%

61	17.292	2412	2415	2421	rVB8	15889	25654	0.61%	0.057%
62	17.416	2433	2436	2441	rVB4	10758	14909	0.36%	0.033%
63	17.475	2443	2446	2448	rBV3	8039	11685	0.28%	0.026%
64	17.616	2467	2470	2473	rBV2	32401	36370	0.87%	0.080%
65	17.786	2495	2499	2506	rVB2	61971	78836	1.88%	0.174%

66	18.004	2533	2536	2543	rVB4	54917	84294	2.01%	0.186%
67	18.163	2560	2563	2568	rBV4	15573	27326	0.65%	0.060%
68	18.204	2568	2570	2575	rVB4	16265	18231	0.44%	0.040%
69	18.251	2575	2578	2584	rVB3	19035	23100	0.55%	0.051%
70	18.310	2584	2588	2593	rBV	41710	58738	1.40%	0.130%

71	18.427	2603	2608	2612	rBV	65819	86576	2.07%	0.191%
72	18.722	2655	2658	2662	rBV6	13066	17965	0.43%	0.040%
73	18.775	2665	2667	2671	rVB4	16620	20542	0.49%	0.045%
74	18.892	2684	2687	2692	rBV4	20759	32277	0.77%	0.071%
75	18.945	2692	2696	2702	rBV2	68074	123104	2.94%	0.272%

76	19.163	2728	2733	2738	rVB	264660	379021	9.05%	0.836%
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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

77	19.292	2752	2755	2756	rBV2	22767	24400	0.58%	0.054%
78	19.380	2767	2770	2777	rVB4	23743	40490	0.97%	0.089%
79	19.498	2784	2790	2803	rBV	2696200	4186989	100.00%	9.237%
80	19.704	2823	2825	2827	rBV	18266	17311	0.41%	0.038%
81	20.022	2877	2879	2883	rVB	42444	45282	1.08%	0.100%
82	20.063	2883	2886	2890	rBV	36546	41230	0.98%	0.091%
83	20.180	2903	2906	2911	rVB2	35519	49095	1.17%	0.108%
84	20.322	2927	2930	2933	rVB2	21605	26098	0.62%	0.058%
85	20.516	2959	2963	2969	rBV	2104551	2268116	54.17%	5.003%
86	20.980	3040	3042	3045	rBV2	27610	26399	0.63%	0.058%
87	21.027	3046	3050	3054	rVV2	122209	169092	4.04%	0.373%
88	21.221	3079	3083	3088	rVB	339945	389880	9.31%	0.860%
89	21.292	3089	3095	3100	rVV2	1626740	2297058	54.86%	5.067%
90	21.333	3100	3102	3109	rVB	172883	201751	4.82%	0.445%
91	21.827	3182	3186	3194	rVB5	122810	236905	5.66%	0.523%
92	21.904	3196	3199	3208	rVB2	101309	145799	3.48%	0.322%
93	22.045	3220	3223	3227	rBV4	64057	88282	2.11%	0.195%
94	22.163	3239	3243	3251	rVB	190893	277358	6.62%	0.612%
95	22.880	3359	3365	3376	rVB3	305253	733360	17.52%	1.618%
96	23.257	3426	3429	3436	rVB3	80212	137003	3.27%	0.302%
97	23.404	3447	3454	3459	rBV	1938346	3563051	85.10%	7.860%
98	23.545	3471	3478	3493	rVB2	1105994	2126019	50.78%	4.690%
99	24.104	3569	3573	3581	rVB	160177	302957	7.24%	0.668%
100	24.857	3696	3701	3713	rVB2	120841	293884	7.02%	0.648%

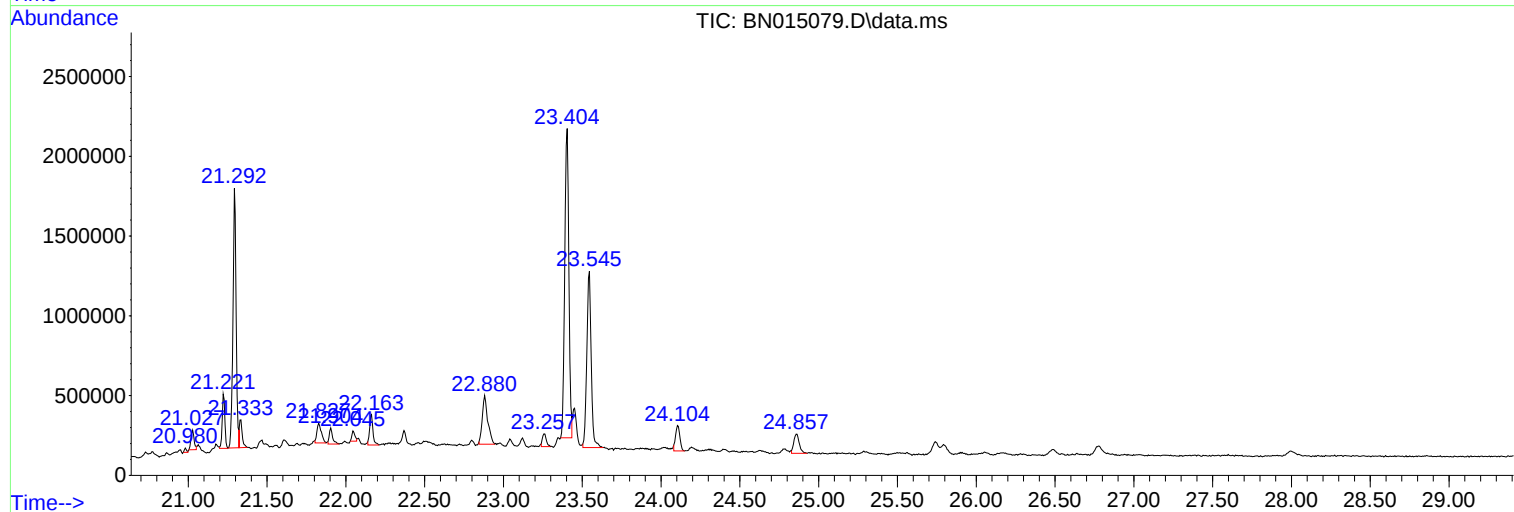
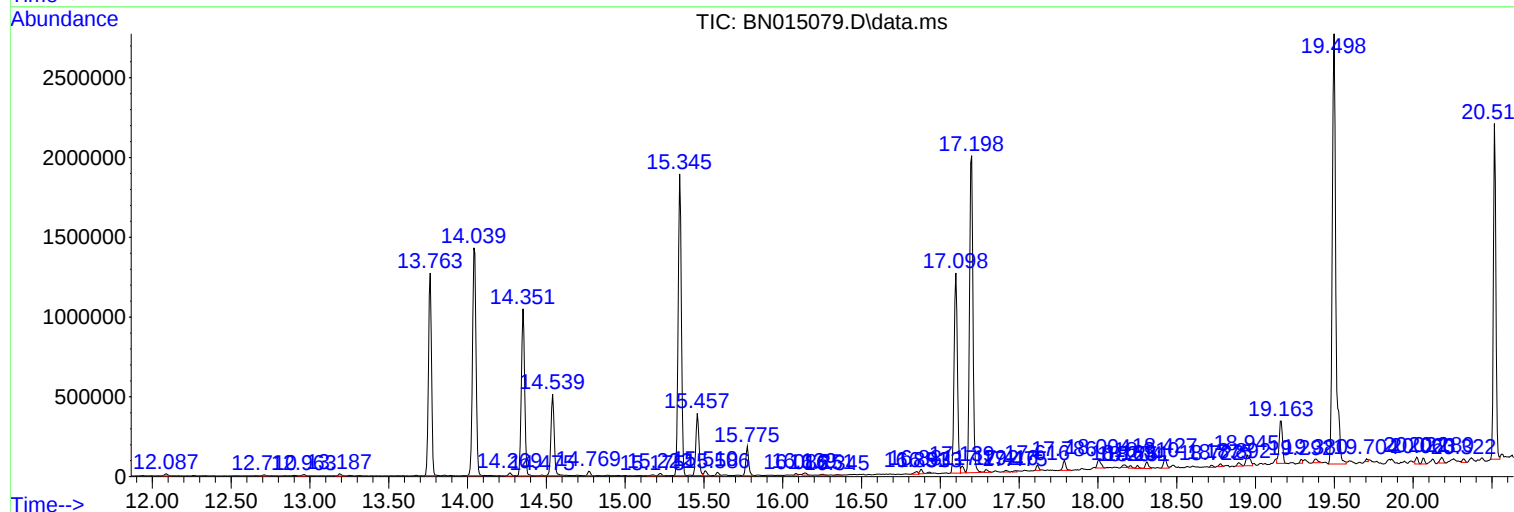
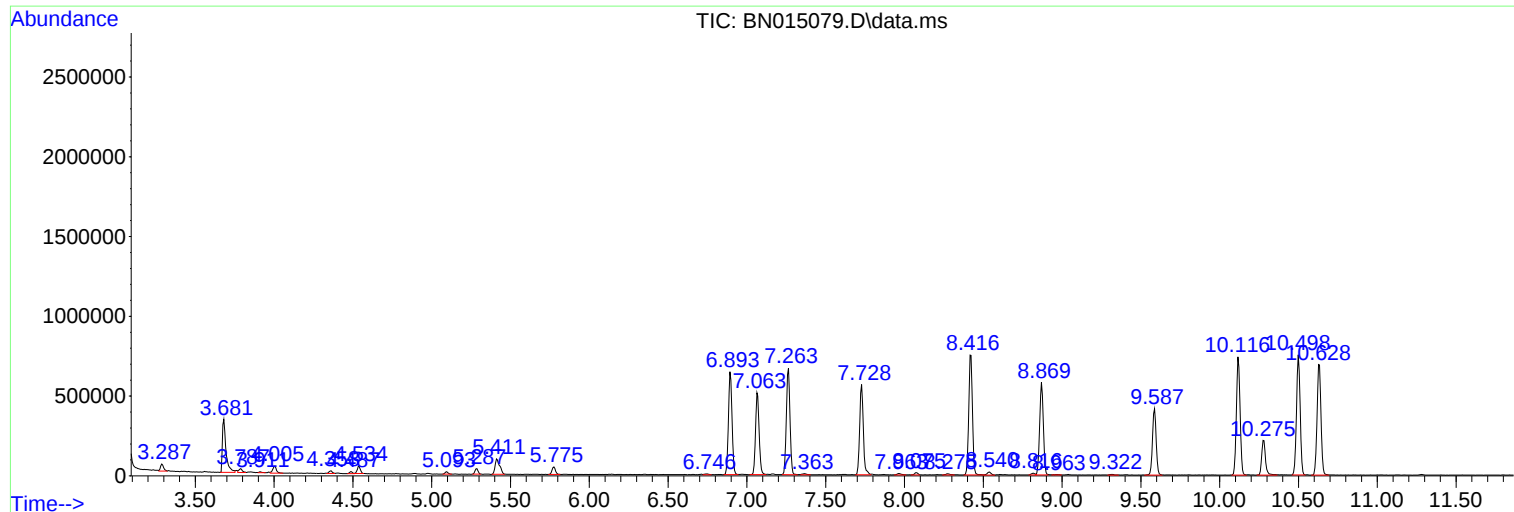
Sum of corrected areas: 45330732

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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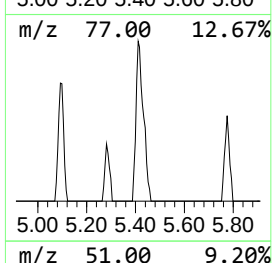
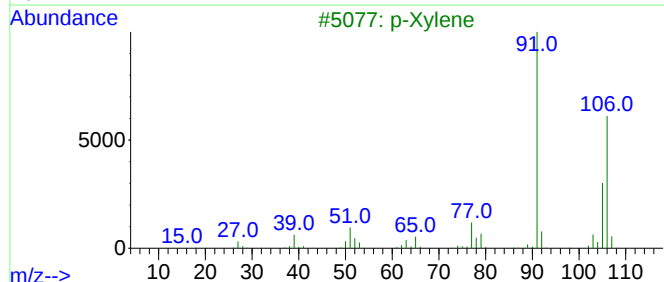
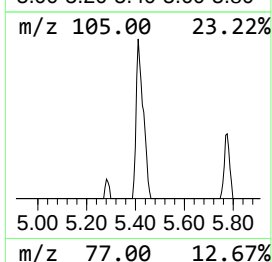
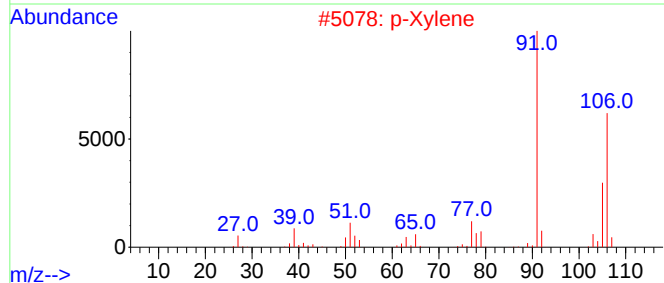
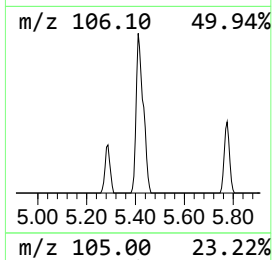
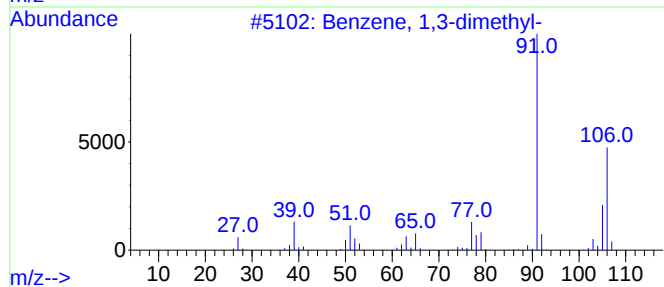
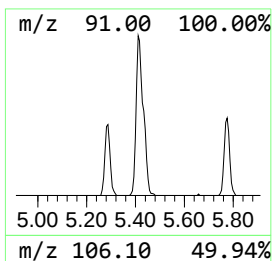
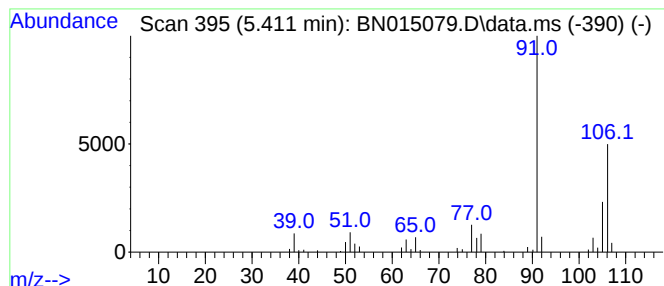
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	4.30 ng/ul	195630	1,4-Dichlorobenzene-d4	7.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			p-Xylene	106	C8H10	000106-42-3	97
3			p-Xylene	106	C8H10	000106-42-3	97
4			o-Xylene	106	C8H10	000095-47-6	95
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95



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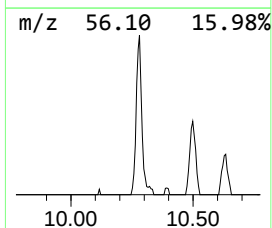
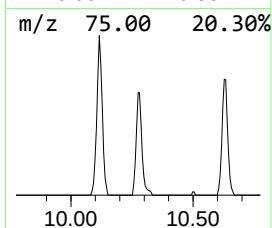
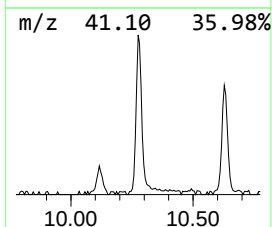
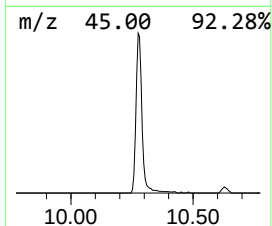
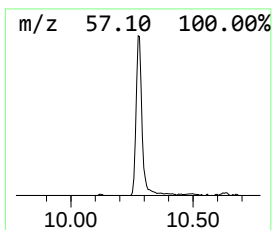
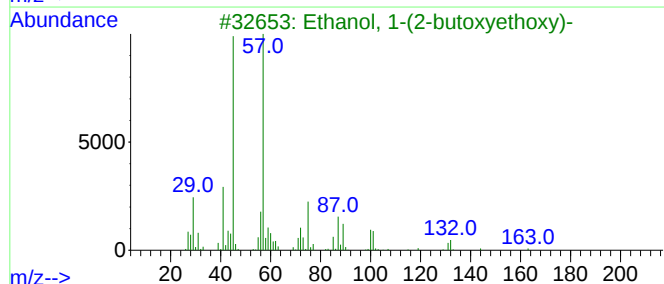
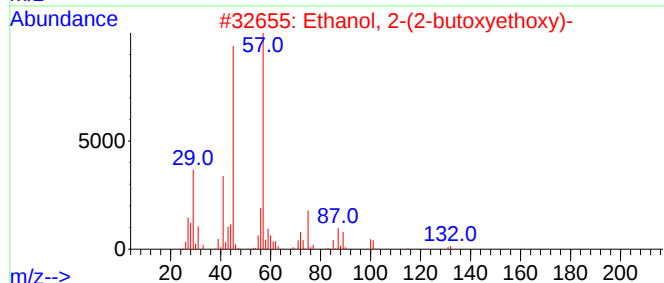
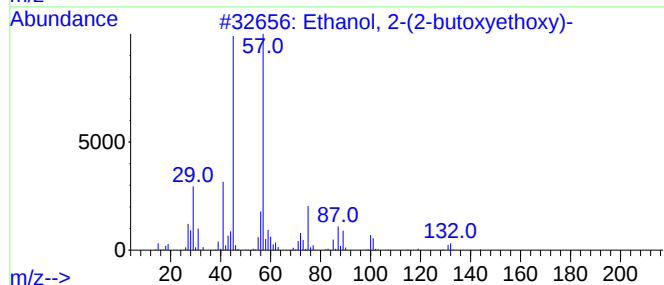
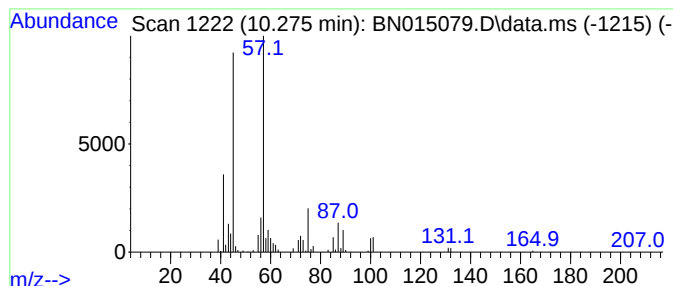
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	5.98 ng/ul	364279	Naphthalene-d8	10.499

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, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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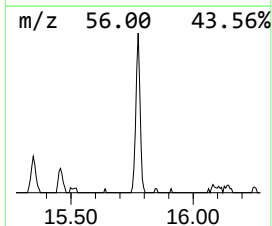
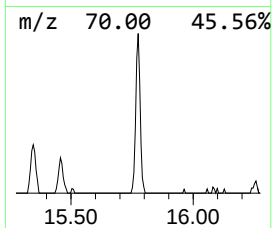
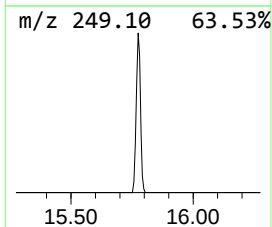
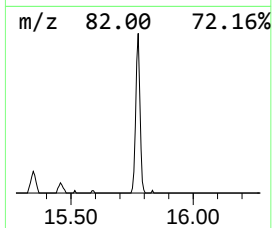
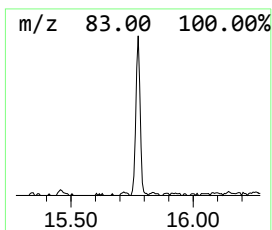
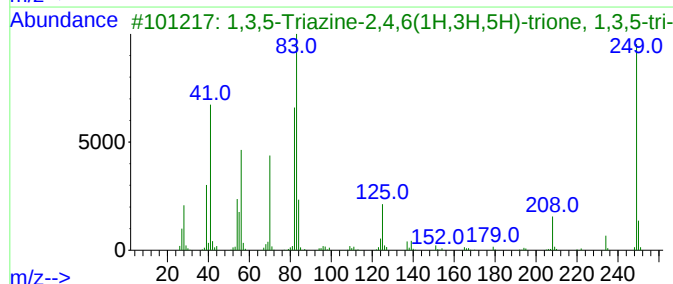
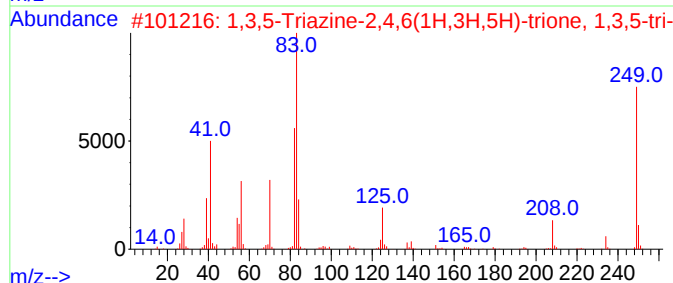
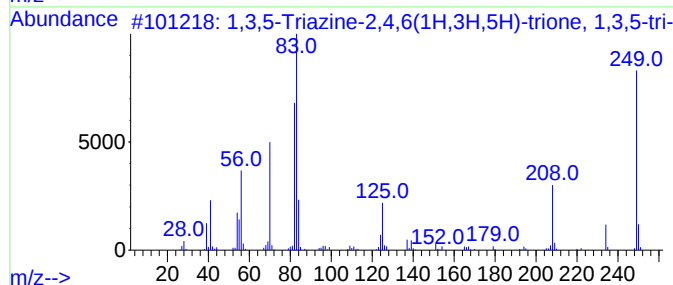
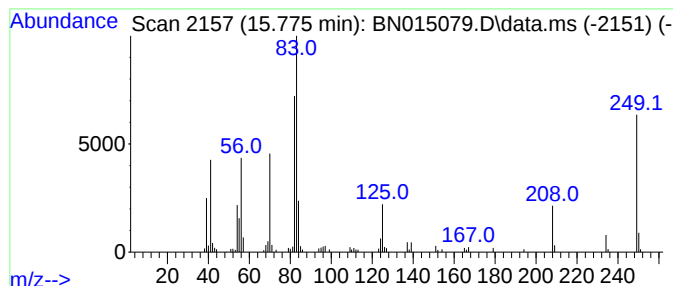
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.63 ng/ul	234910	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98		
2	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	96		
3	1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	96		
4	N-Carboxy-1-[4-chlorophenyl]-2-[...	295	C15H18ClNO3	058158-38-6	37		
5	Cyclohexane, 1,1'-(1,4-butanediyl...	222	C16H30	006165-44-2	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015079.D
Acq On : 25 Jun 2021 15:55
Operator : CG/JU
Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD34

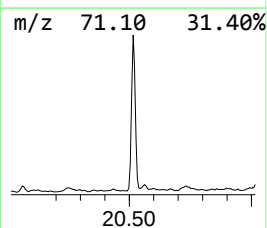
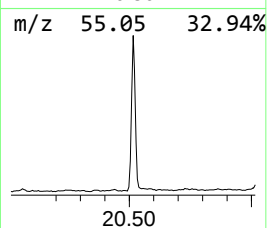
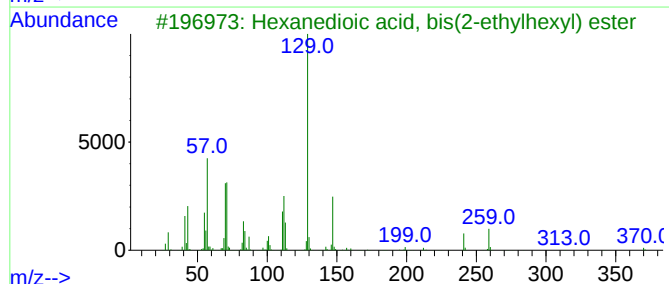
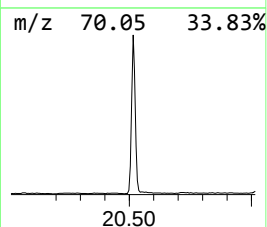
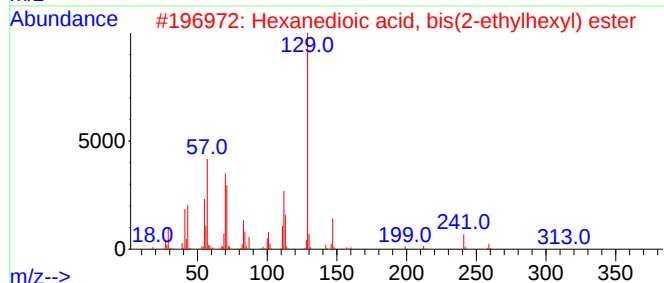
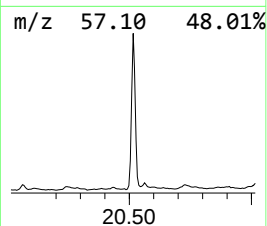
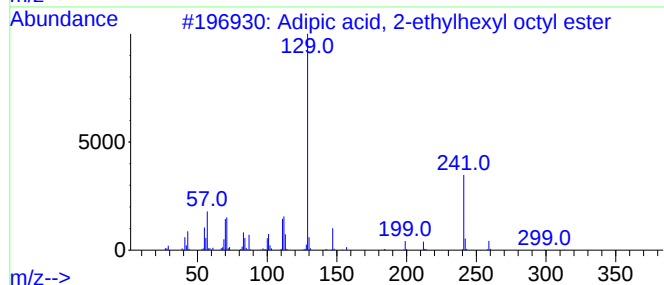
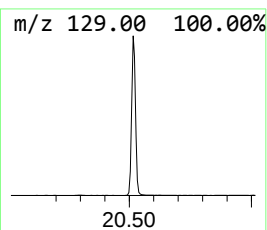
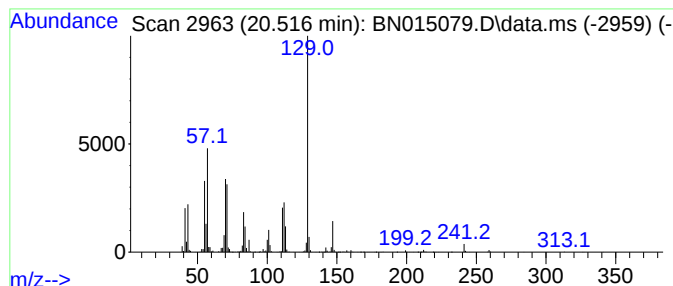
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Adipic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	19.75 ng/ul	2268120	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	91
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015079.D
Acq On : 25 Jun 2021 15:55
Operator : CG/JU
Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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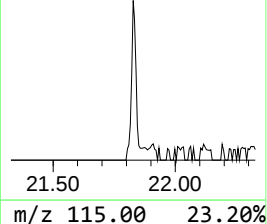
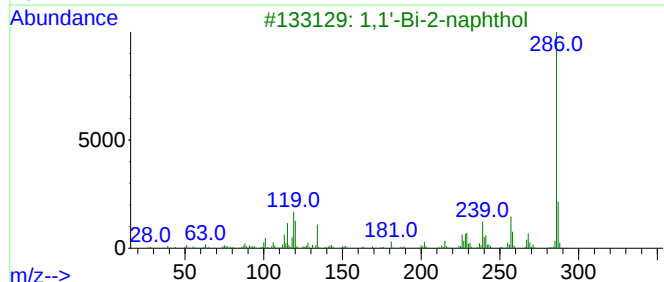
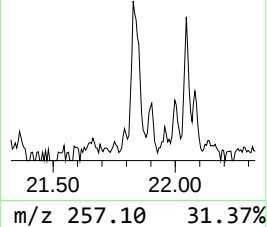
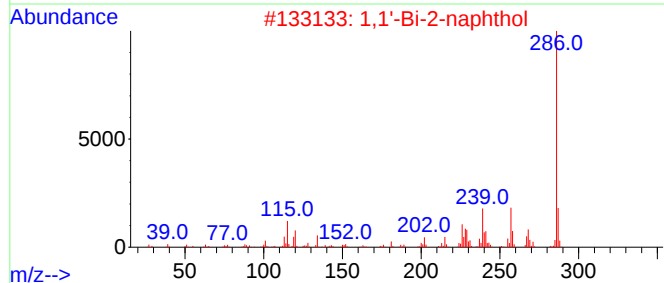
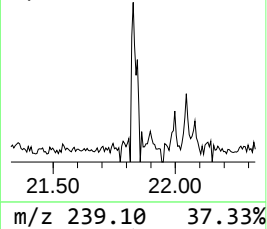
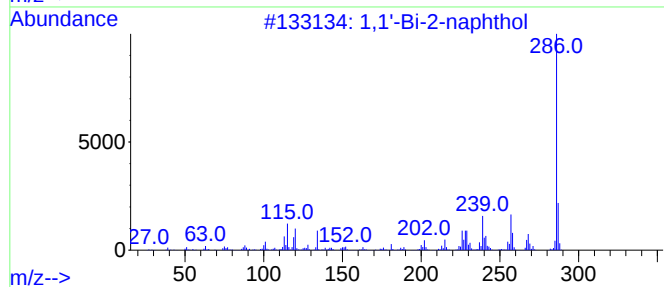
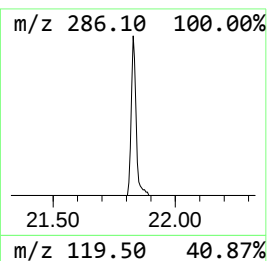
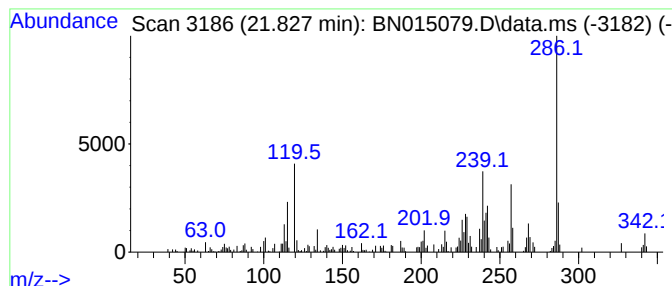
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Bi-2-naphthol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.827	2.06 ng/ul	236905	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Bi-2-naphthol	286	C20H14O2	000602-09-5	98		
2	1,1'-Bi-2-naphthol	286	C20H14O2	000602-09-5	96		
3	1,1'-Bi-2-naphthol	286	C20H14O2	000602-09-5	86		
4	5-(Acridin-9-yl)-2-aminophenol	286	C19H14N2O	1000326-85-0	64		
5	[1,1'-Binaphthalene]-4,4'-diol	286	C20H14O2	001446-34-0	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015079.D
Acq On : 25 Jun 2021 15:55
Operator : CG/JU
Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

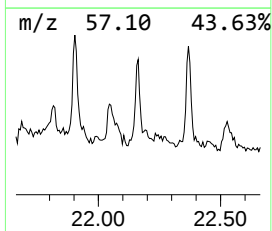
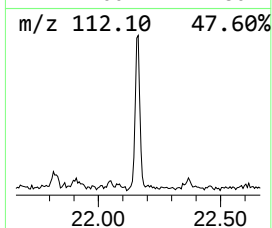
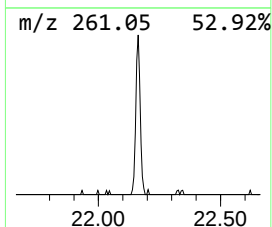
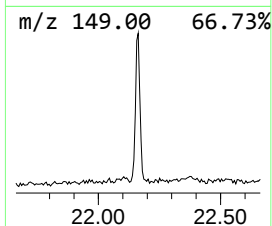
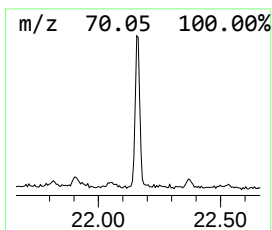
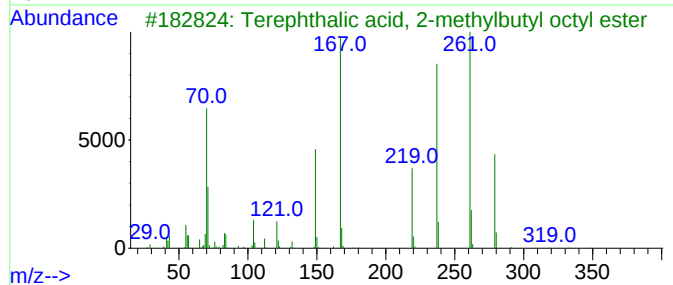
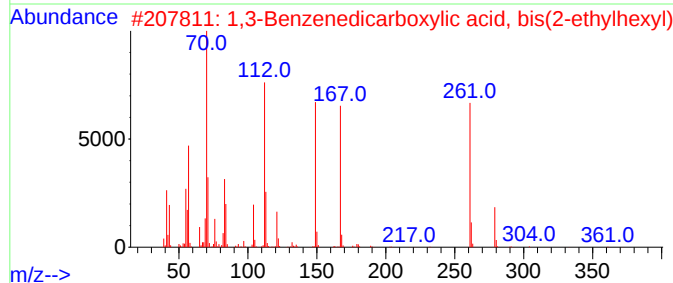
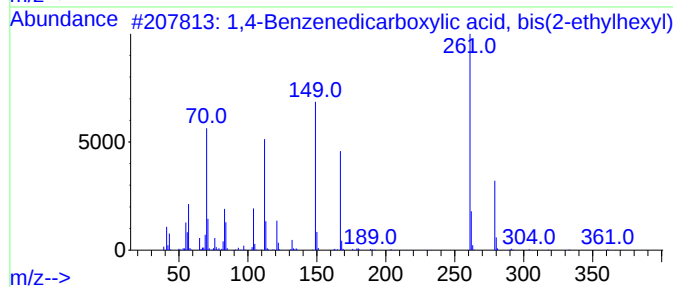
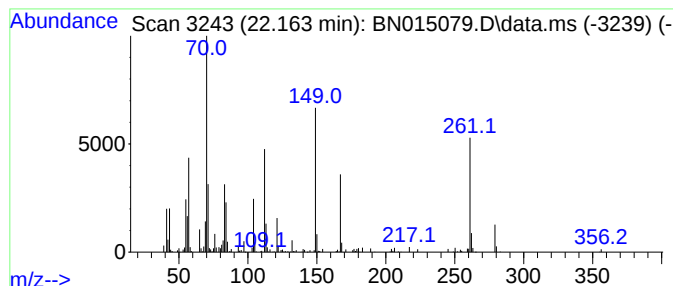
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,4-Benzenedicarboxylic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.163	2.41 ng/ul	277358	Chrysene-d12	21.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	91
2	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	52
3	Terephthalic acid, 2-methylbutyl...	348	C21H32O4	1000323-90-4	50
4	Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	49
5	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015079.D
Acq On : 25 Jun 2021 15:55
Operator : CG/JU
Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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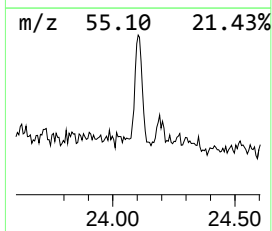
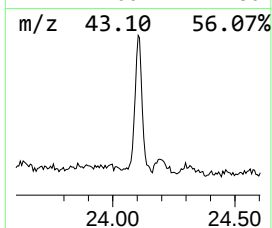
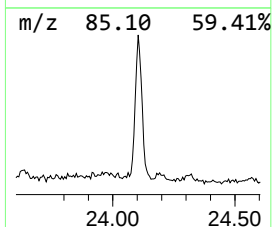
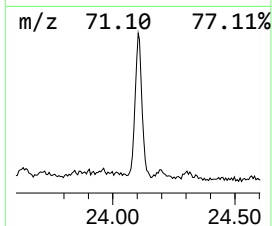
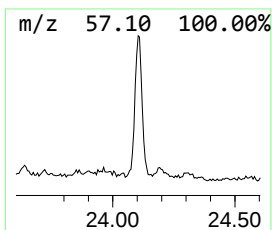
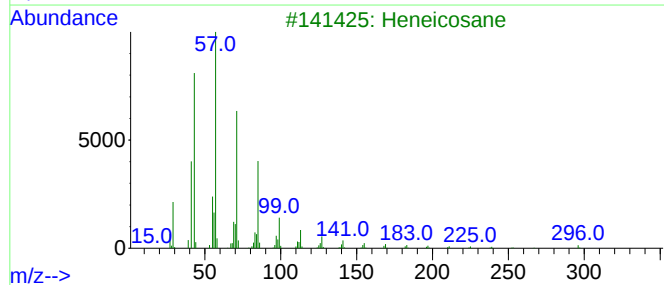
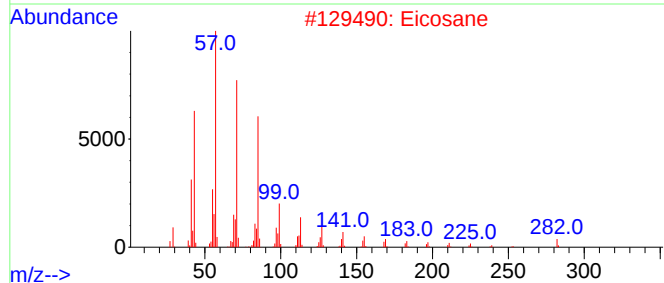
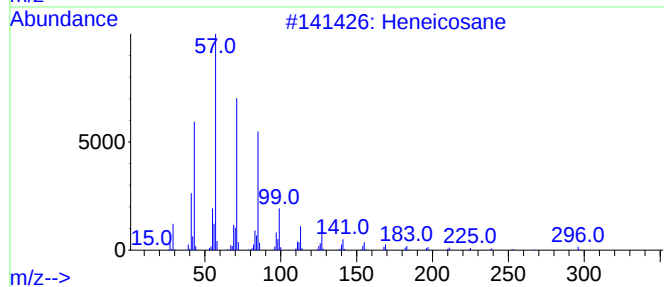
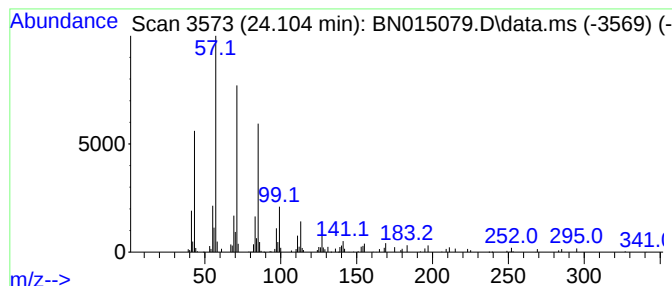
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.104	2.85 ng/ul	302957	Perylene-d12	23.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Hexacosane	366	C26H54	000630-01-3	87
5		Heptadecane	240	C17H36	000629-78-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015079.D
Acq On : 25 Jun 2021 15:55
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Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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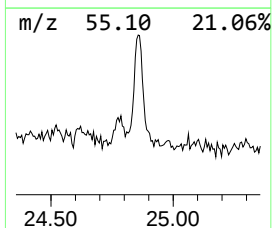
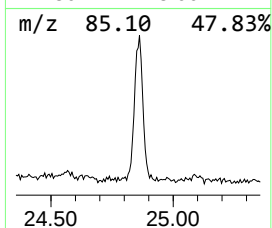
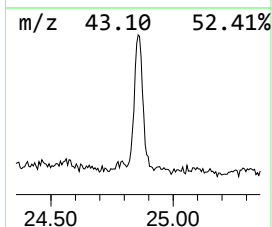
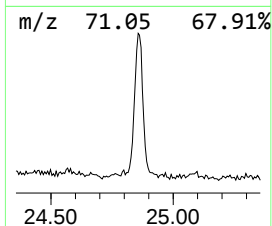
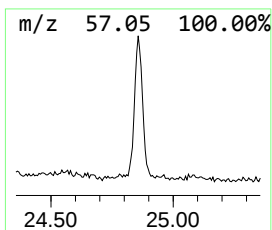
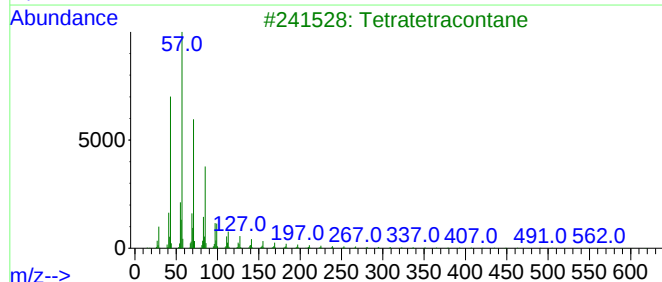
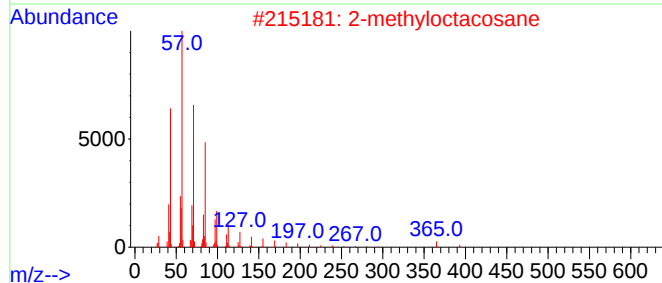
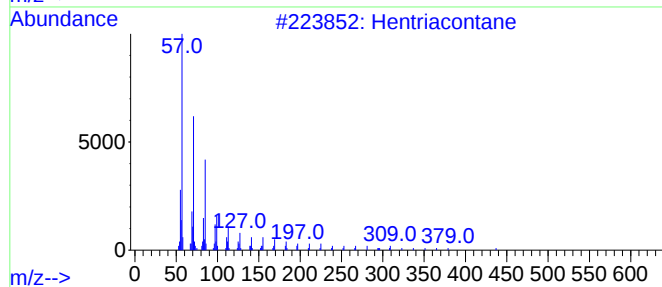
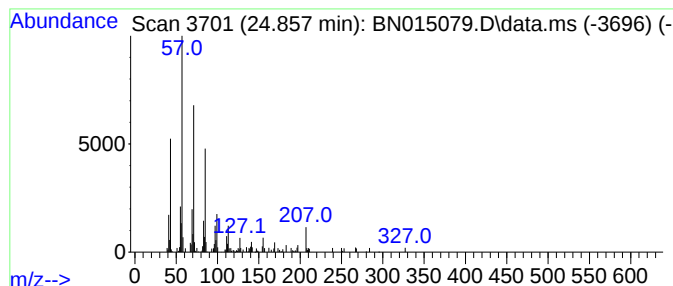
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.857	2.76 ng/ul	293884	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			2-methyloctacosane	408	C29H60	1000376-72-8	87
3			Tetratetracontane	619	C44H90	007098-22-8	87
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Operator : CG/JU
Sample : M2680-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Ethanol, 2-(2-b...	10.275	6.0	ng/ul	364279	2	10.499	1217440	20.0
1,3,5-Triazine-...	15.775	2.6	ng/ul	234910	4	17.098	1786890	20.0
Adipic acid, 2-...	20.516	19.8	ng/ul	2268120	5	21.292	2297060	20.0
1,1'-Bi-2-naphthol	21.827	2.1	ng/ul	236905	5	21.292	2297060	20.0
1,4-Benzenedica...	22.163	2.4	ng/ul	277358	5	21.292	2297060	20.0
(DEL) Alkane: S...	24.104	2.9	ng/ul	302957	6	23.545	2126020	20.0
(DEL) Alkane: S...	24.857	2.8	ng/ul	293884	6	23.545	2126020	20.0