

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
 Data File : BN015041.D
 Acq On : 24 Jun 2021 15:30
 Operator : CG/JU
 Sample : M2682-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGDB0

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

Signal : TIC: BN015041.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	43	rVB2	13255	22875	0.67%	0.079%
2	3.687	99	102	113	rBV	53370	88892	2.60%	0.305%
3	3.787	115	119	124	rVB2	17694	25244	0.74%	0.087%
4	3.911	136	140	147	rVB3	8307	15986	0.47%	0.055%
5	3.999	151	155	163	rVB	47460	68100	1.99%	0.234%
6	4.211	187	191	198	rVB	16141	22520	0.66%	0.077%
7	4.358	212	216	221	rVB5	9824	15056	0.44%	0.052%
8	4.487	235	238	242	rVB2	6936	9441	0.28%	0.032%
9	4.534	242	246	253	rVV	26758	38325	1.12%	0.132%
10	4.611	255	259	264	rVB3	7424	10625	0.31%	0.036%
11	4.887	302	306	311	rVB2	8633	13092	0.38%	0.045%
12	5.287	369	374	382	rVB	19579	29827	0.87%	0.102%
13	5.411	390	395	404	rVB	57137	114128	3.34%	0.392%
14	5.775	452	457	462	rVB	27382	40366	1.18%	0.139%
15	6.893	639	647	656	rVB	176698	286364	8.39%	0.983%
16	7.064	671	676	684	rVB	140634	220707	6.46%	0.758%
17	7.264	704	710	718	rBV	180396	290455	8.51%	0.997%
18	7.364	724	727	735	rBV4	4396	7380	0.22%	0.025%
19	7.728	781	789	796	rVB	637185	1002605	29.36%	3.442%
20	8.416	899	906	913	rBV	211745	321280	9.41%	1.103%
21	8.534	921	926	934	rVB	20932	33449	0.98%	0.115%
22	8.816	968	974	977	rBV2	6234	8212	0.24%	0.028%
23	8.869	977	983	990	rVB	142416	228576	6.69%	0.785%
24	9.587	1098	1105	1112	rVB	91945	146428	4.29%	0.503%
25	10.116	1189	1195	1203	rVB	186311	298226	8.73%	1.024%
26	10.281	1218	1223	1228	rBV2	8724	13224	0.39%	0.045%
27	10.499	1253	1260	1267	rBV	849191	1373875	40.24%	4.717%
28	10.628	1276	1282	1291	rBV	161040	273276	8.00%	0.938%
29	11.275	1388	1392	1395	rBV4	4673	6396	0.19%	0.022%
30	13.193	1712	1718	1721	rBV2	5488	7267	0.21%	0.025%
31	13.493	1764	1769	1772	rBV2	6121	5828	0.17%	0.020%
32	13.681	1796	1801	1808	rVB	85119	124357	3.64%	0.427%
33	13.763	1808	1815	1820	rBV	339247	467750	13.70%	1.606%
34	14.040	1856	1862	1865	rBV	366956	559164	16.38%	1.920%
35	14.075	1865	1868	1875	rVB	288566	385374	11.29%	1.323%
36	14.204	1885	1890	1895	rBV2	26384	33928	0.99%	0.116%

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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
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37	14.269	1897	1901	1906	rVV	9940	12268	0.36%	0.042%
38	14.351	1908	1915	1923	rVV	1214520	1717824	50.31%	5.898%
39	14.428	1925	1928	1932	rVB5	6927	9122	0.27%	0.031%
40	14.534	1942	1946	1953	rBV2	15546	23347	0.68%	0.080%
41	15.222	2059	2063	2067	rBV3	7034	9029	0.26%	0.031%
42	15.345	2079	2084	2090	rVV	514289	741934	21.73%	2.547%
43	15.457	2097	2103	2106	rBV5	12111	17180	0.50%	0.059%
44	15.522	2111	2114	2118	rVB2	7852	11605	0.34%	0.040%
45	15.775	2150	2157	2163	rBV2	79447	108539	3.18%	0.373%
46	15.851	2165	2170	2174	rVB5	3031	5825	0.17%	0.020%
47	16.040	2194	2202	2207	rBV	32061	46333	1.36%	0.159%
48	16.087	2207	2210	2215	rVB5	5122	8118	0.24%	0.028%
49	16.604	2292	2298	2302	rBV6	3787	7919	0.23%	0.027%
50	16.763	2319	2325	2330	rBV8	7452	11877	0.35%	0.041%
51	16.881	2339	2345	2348	rBV3	16271	21427	0.63%	0.074%
52	16.934	2350	2354	2362	rVB3	15127	23291	0.68%	0.080%
53	17.098	2376	2382	2387	rVV	1370596	1959766	57.40%	6.729%
54	17.139	2387	2389	2393	rVV	53041	81500	2.39%	0.280%
55	17.192	2393	2398	2410	rVV2	560815	864045	25.31%	2.967%
56	17.304	2412	2417	2421	rVV8	7945	16686	0.49%	0.057%
57	17.339	2421	2423	2429	rVV5	4980	8072	0.24%	0.028%
58	17.475	2441	2446	2450	rBV2	14061	22815	0.67%	0.078%
59	17.616	2466	2470	2474	rBV2	14242	20777	0.61%	0.071%
60	17.786	2497	2499	2504	rVB5	6969	10695	0.31%	0.037%
61	17.969	2528	2530	2533	rBV4	7791	11174	0.33%	0.038%
62	18.004	2533	2536	2544	rVB4	51428	80177	2.35%	0.275%
63	18.169	2559	2564	2568	rBV6	11286	19029	0.56%	0.065%
64	18.251	2576	2578	2582	rVB4	8124	8899	0.26%	0.031%
65	18.310	2585	2588	2592	rBV	19177	24798	0.73%	0.085%
66	18.428	2603	2608	2612	rBV	95027	127443	3.73%	0.438%
67	18.781	2665	2668	2670	rBV4	7734	10850	0.32%	0.037%
68	18.892	2684	2687	2692	rBV6	7707	12038	0.35%	0.041%
69	18.951	2693	2697	2704	rBV	28689	54025	1.58%	0.185%
70	19.086	2714	2720	2725	rBV	2110131	2447355	71.68%	8.403%
71	19.163	2728	2733	2738	rVB	98534	138441	4.05%	0.475%
72	19.292	2752	2755	2758	rBV4	31606	42686	1.25%	0.147%
73	19.498	2784	2790	2794	rBV2	740529	1079133	31.60%	3.705%
74	19.898	2854	2858	2867	rVB	330993	389356	11.40%	1.337%
75	20.069	2883	2887	2891	rVB	31079	34296	1.00%	0.118%
76	20.128	2893	2897	2901	rVV4	14648	21623	0.63%	0.074%

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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

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77	20.180	2904	2906	2911	rVB6	10241	11966	0.35%	0.041%
78	20.522	2959	2964	2969	rBV	3015355	3414450	100.00%	11.723%
79	20.563	2969	2971	2976	rVB	37640	44111	1.29%	0.151%
80	20.863	3020	3022	3025	rVB4	11435	11183	0.33%	0.038%
81	20.951	3034	3037	3044	rVB5	46548	72796	2.13%	0.250%
82	21.027	3046	3050	3055	rBV	101024	131262	3.84%	0.451%
83	21.157	3068	3072	3074	rBV3	38153	51833	1.52%	0.178%
84	21.227	3080	3084	3088	rVB	126390	140390	4.11%	0.482%
85	21.292	3089	3095	3100	rBV	1571389	2158265	63.21%	7.410%
86	21.469	3122	3125	3137	rVB2	51783	85141	2.49%	0.292%
87	21.910	3195	3200	3209	rBV3	187402	425904	12.47%	1.462%
88	22.069	3223	3227	3232	rVB2	117704	164607	4.82%	0.565%
89	22.163	3239	3243	3246	rBV	87730	119774	3.51%	0.411%
90	22.374	3275	3279	3288	rVB	111476	182411	5.34%	0.626%
91	22.798	3349	3351	3356	rVB5	48768	60844	1.78%	0.209%
92	22.886	3360	3366	3371	rBV	318697	512945	15.02%	1.761%
93	23.398	3447	3453	3458	rBV	428715	798140	23.38%	2.740%
94	23.457	3459	3463	3470	rVB2	149615	247337	7.24%	0.849%
95	23.545	3470	3478	3486	rBV2	1141619	2198262	64.38%	7.548%
96	24.110	3568	3574	3583	rVB	178291	352226	10.32%	1.209%
97	24.863	3696	3702	3711	rVB2	116015	265509	7.78%	0.912%
98	25.745	3847	3852	3863	rVB4	64672	161508	4.73%	0.555%
99	26.786	4023	4029	4040	rVB4	64887	196946	5.77%	0.676%
100	27.915	4213	4221	4231	rBV5	131225	447464	13.11%	1.536%

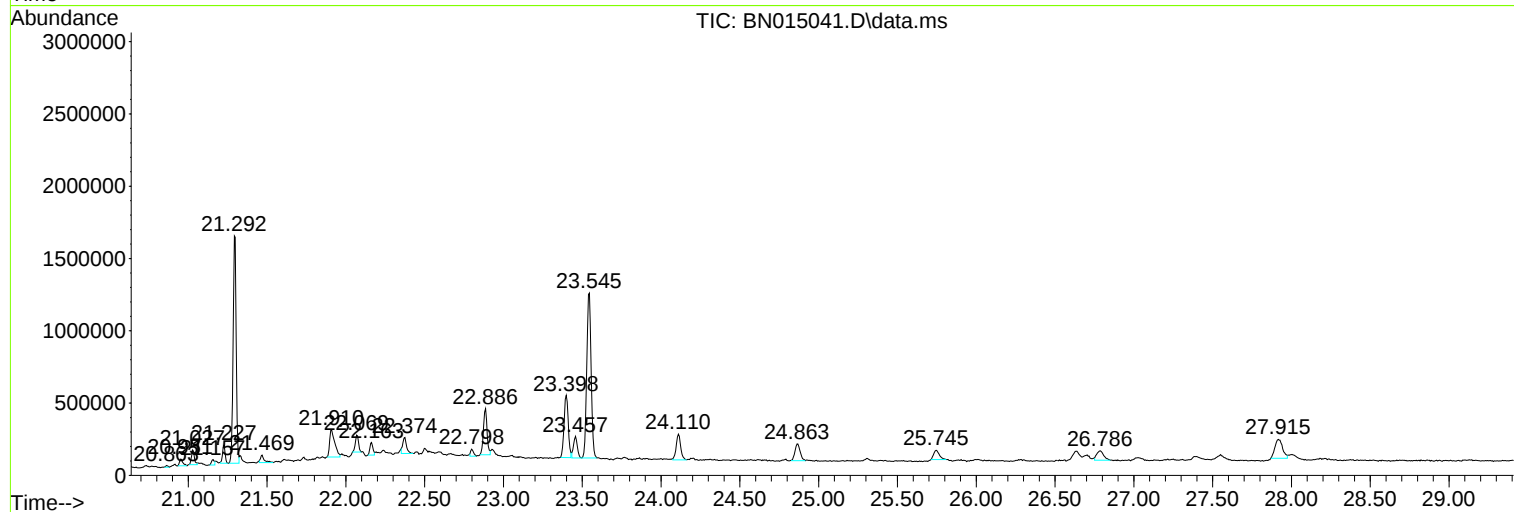
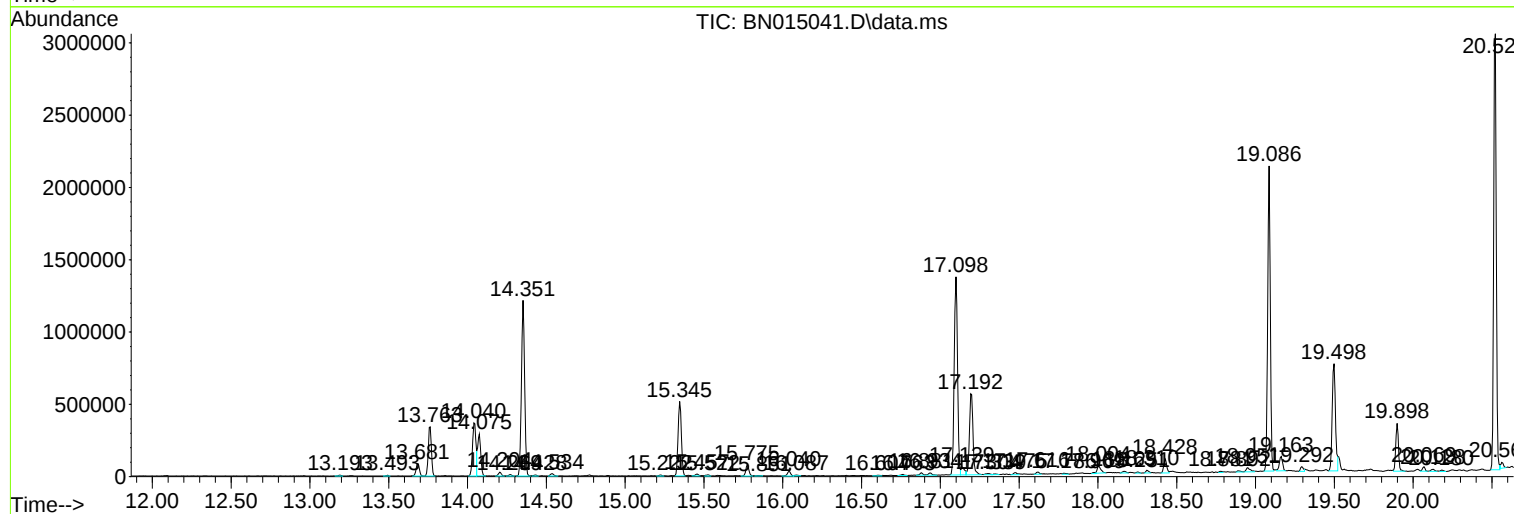
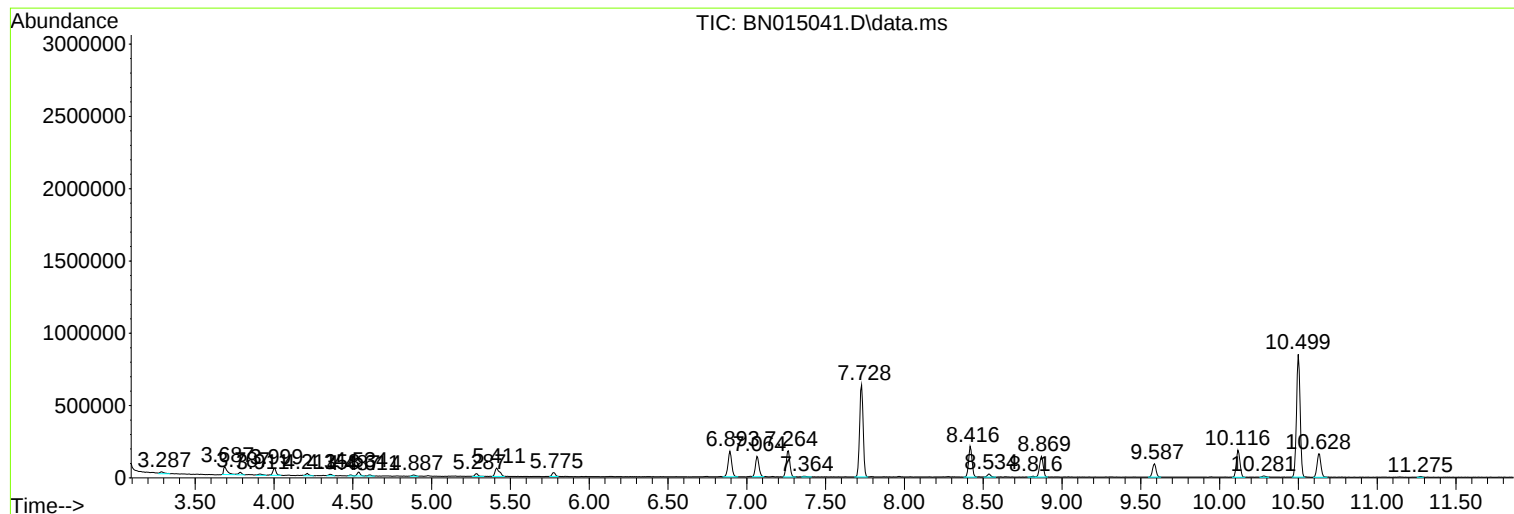
Sum of corrected areas: 29125189

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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



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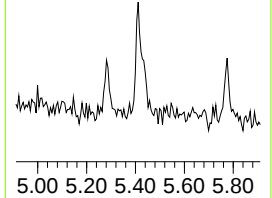
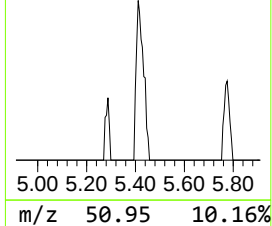
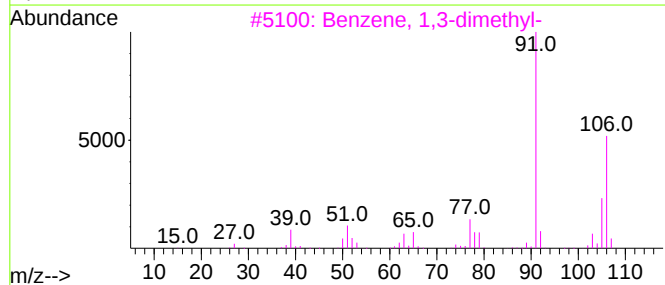
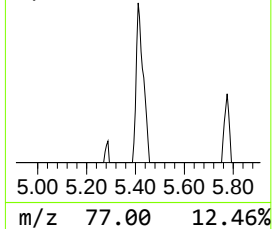
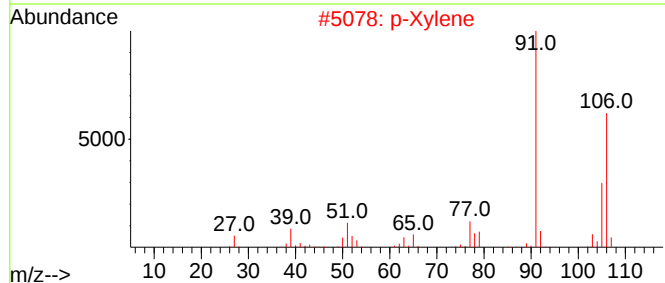
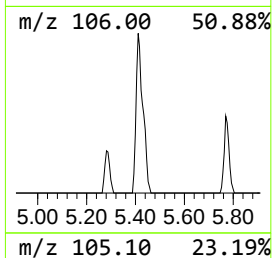
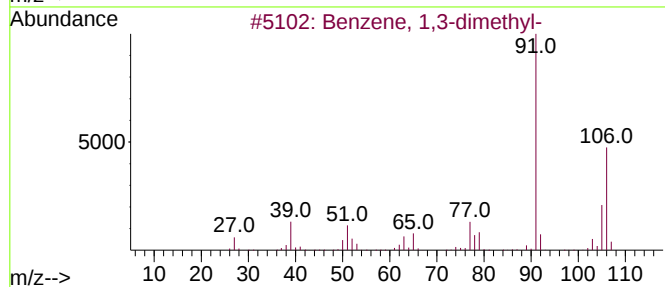
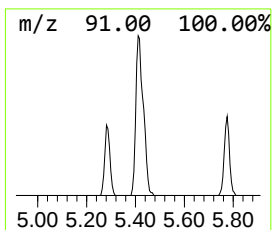
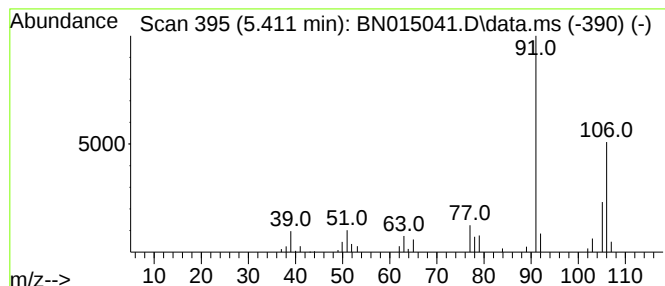
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.411	2.28 ng/ul	114128	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	95
2			p-Xylene	106	C8H10	000106-42-3	95
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
4			p-Xylene	106	C8H10	000106-42-3	94
5			o-Xylene	106	C8H10	000095-47-6	94



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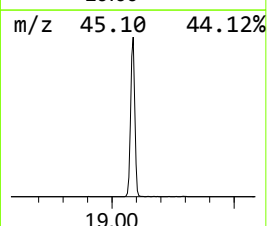
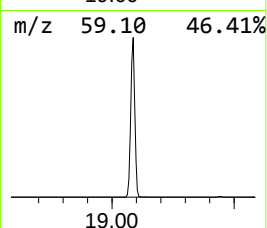
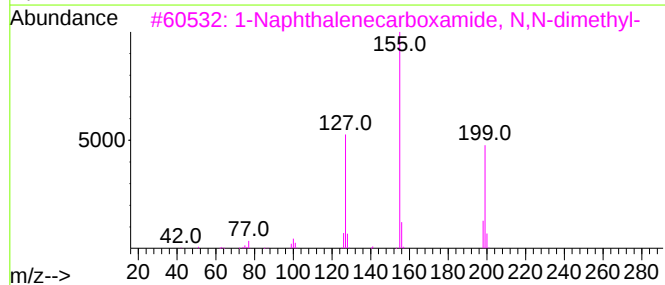
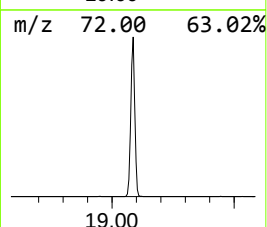
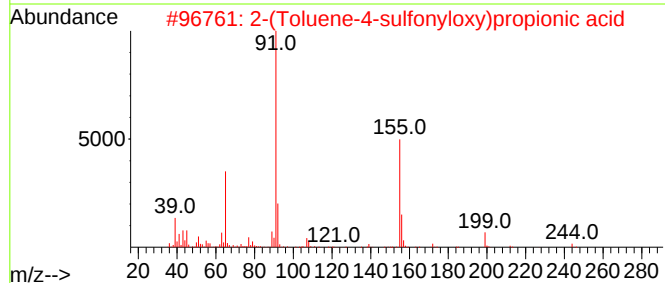
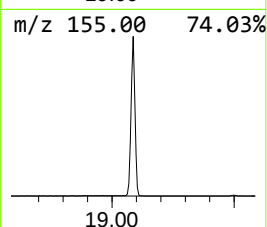
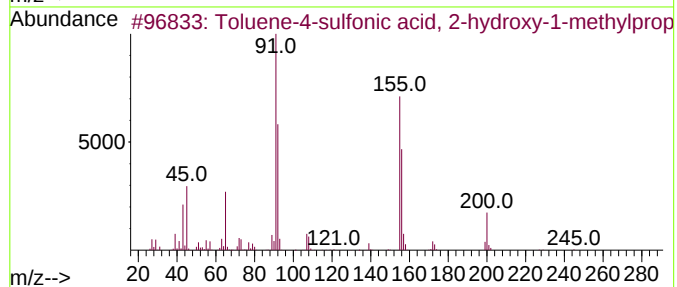
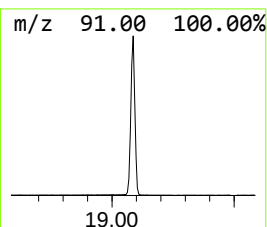
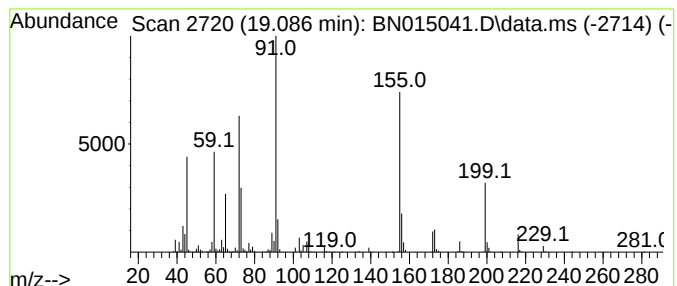
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	24.98 ng/ul	2447360	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene-4-sulfonic acid, 2-hydro...	244	C11H16O4S	1000192-74-7	38
2			2-(Toluene-4-sulfonyloxy)propion...	244	C10H12O5S	1000210-27-2	22
3			1-Naphthalenecarboxamide, N,N-di...	199	C13H13NO	003815-24-5	22
4			Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	18
5			2-Naphthalenecarboxamide, N,N-di...	199	C13H13NO	013577-85-0	16



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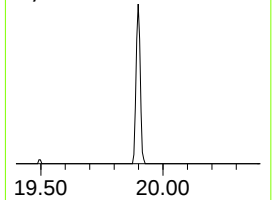
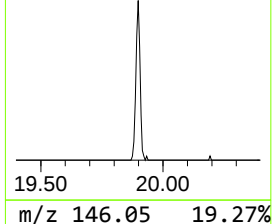
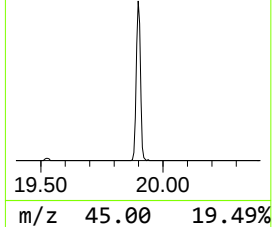
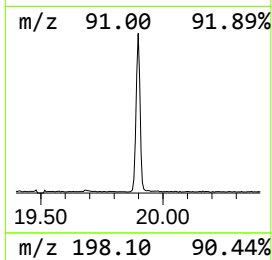
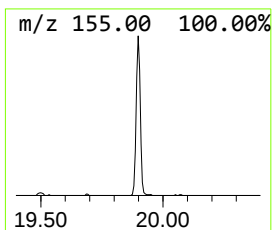
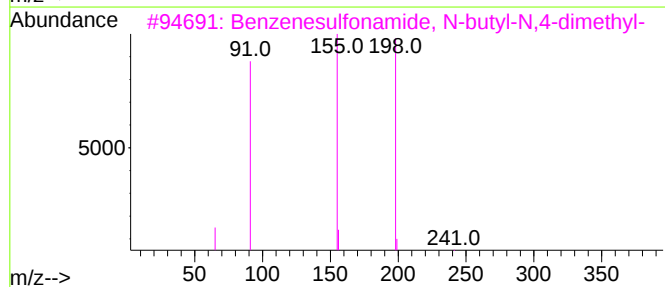
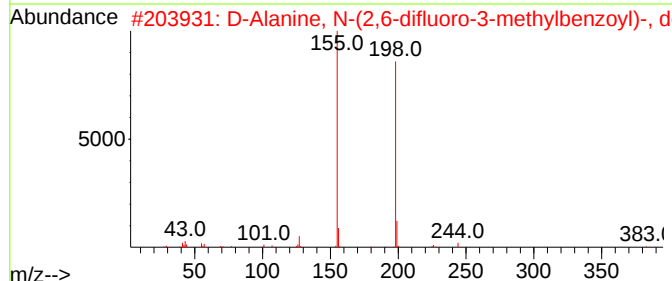
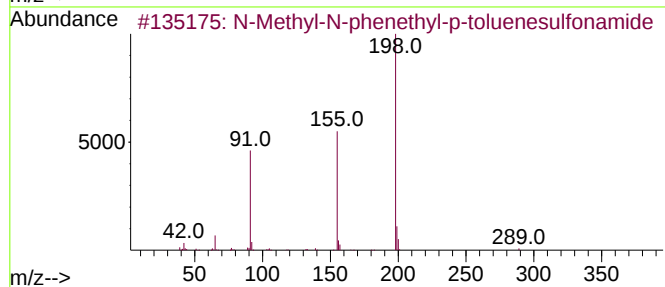
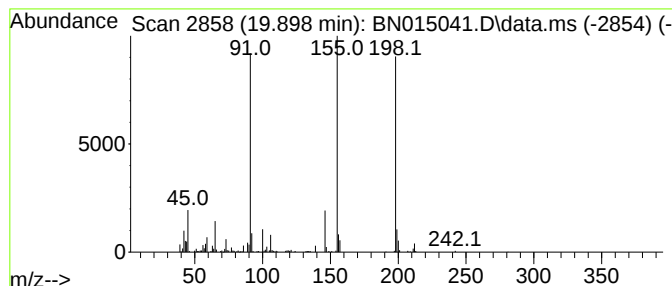
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 N-Methyl-N-phenethyl-p-tolu... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.898	3.61 ng/ul	389356	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Methyl-N-phenethyl-p-toluenesu...	289	C16H19NO2S	025566-59-0	59
2			D-Alanine, N-(2,6-difluoro-3-met...	383	C21H31F2NO3	1000348-39-1	50
3			Benzenesulfonamide, N-butyl-N,4-...	241	C12H19NO2S	010285-91-3	50
4			D-Alanine, N-(2,6-difluoro-3-met...	397	C22H33F2NO3	1000348-39-2	50
5			D-Alanine, N-(2,6-difluoro-3-met...	327	C17H23F2NO3	1000348-38-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
Acq On : 24 Jun 2021 15:30
Operator : CG/JU
Sample : M2682-01
Misc :
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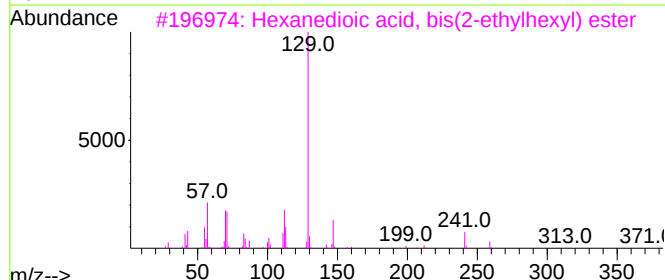
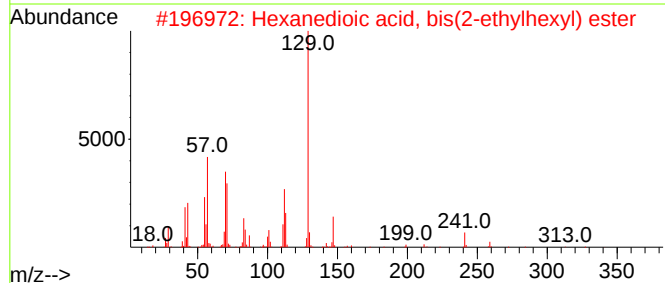
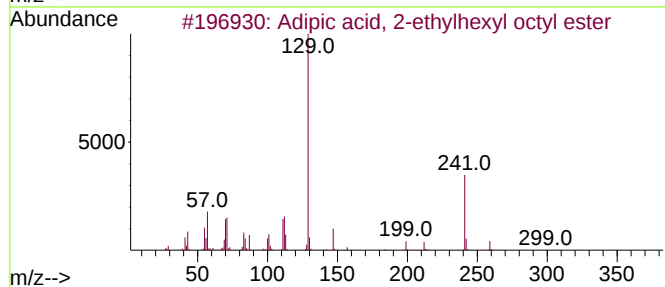
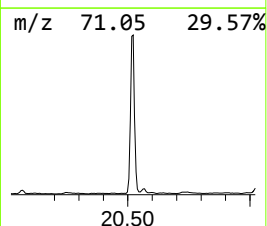
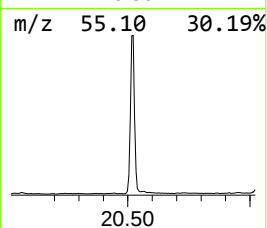
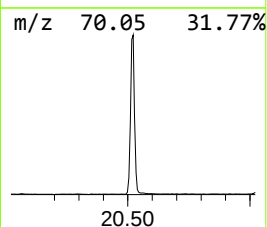
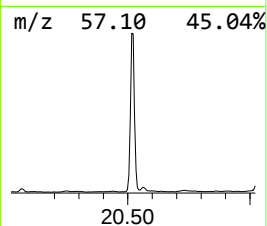
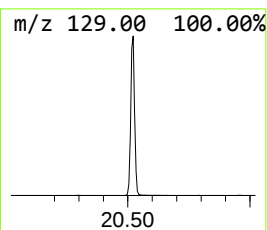
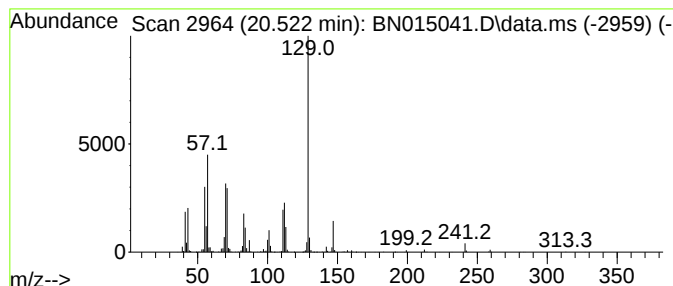
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.522	31.64 ng/ul	3414450	Chrysene-d12	21.298	
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	87
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
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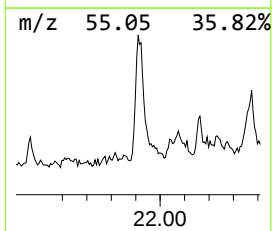
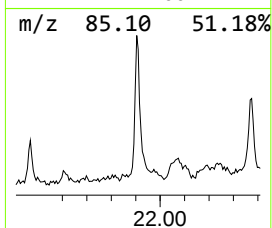
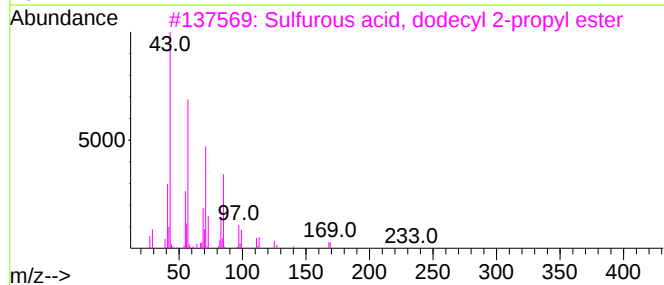
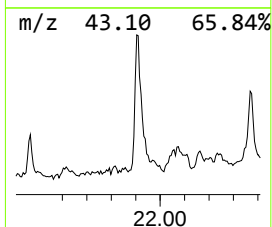
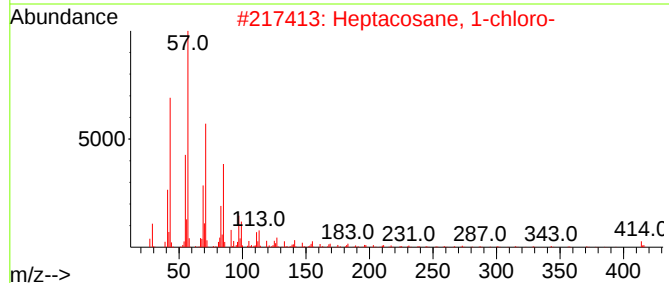
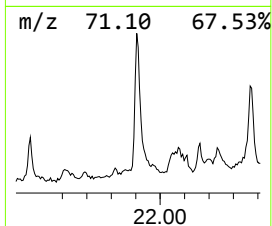
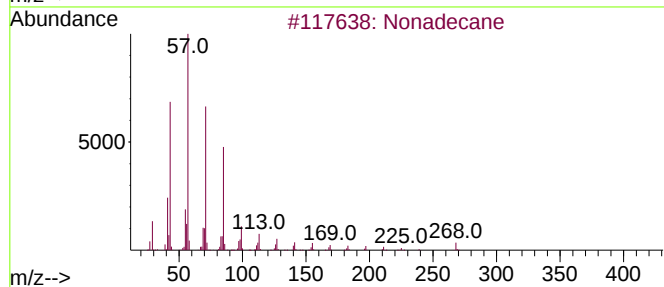
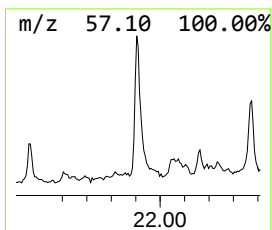
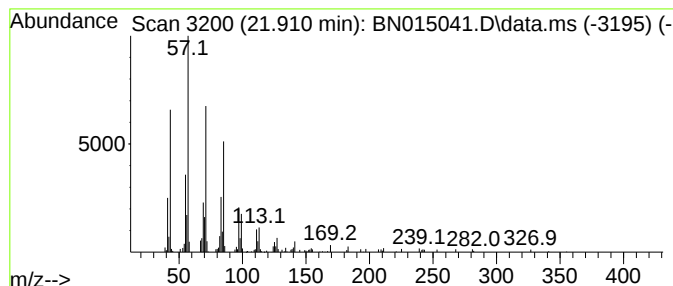
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.910	3.95 ng/ul	425904	Chrysene-d12		21.298	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	91
3	Sulfurous acid, dodecyl 2-propyl...		292	C15H32O3S	1000309-12-3	91
4	Tritetracontane		605	C43H88	007098-21-7	91
5	Sulfurous acid, butyl dodecyl ester		306	C16H34O3S	1000309-17-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
Acq On : 24 Jun 2021 15:30
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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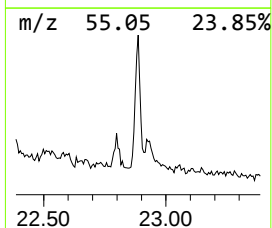
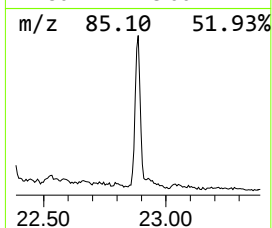
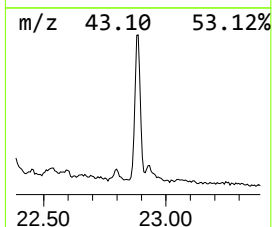
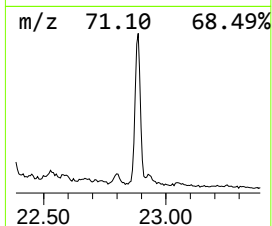
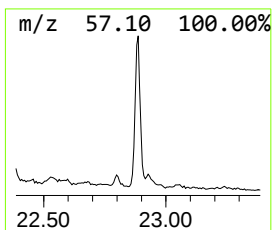
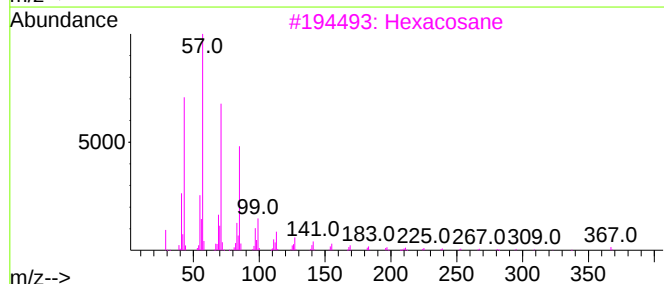
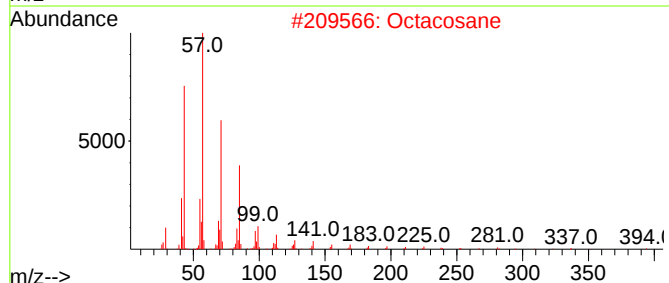
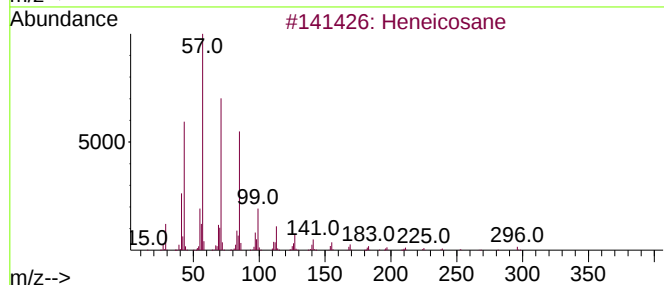
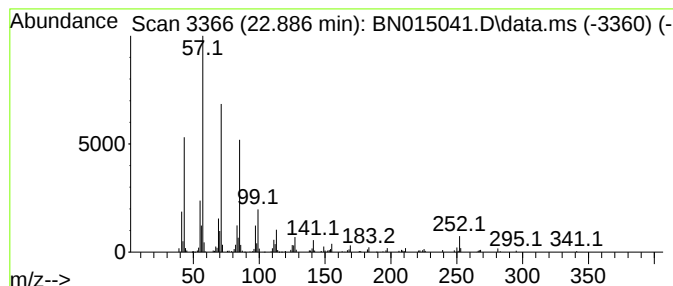
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.886	4.67 ng/ul	512945	Perylene-d12	23.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Octacosane	394	C28H58	000630-02-4	87
3			Hexacosane	366	C26H54	000630-01-3	87
4			2-methyloctacosane	408	C29H60	1000376-72-8	87
5			Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
Acq On : 24 Jun 2021 15:30
Operator : CG/JU
Sample : M2682-01
Misc :
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Instrument :
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ClientSampleId :
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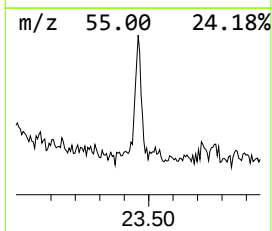
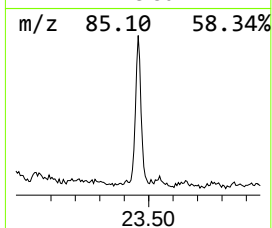
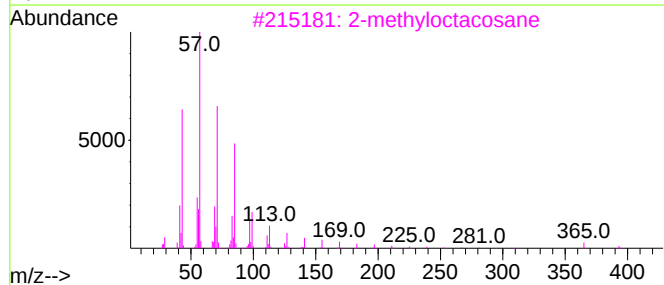
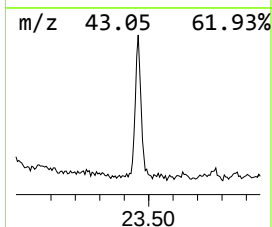
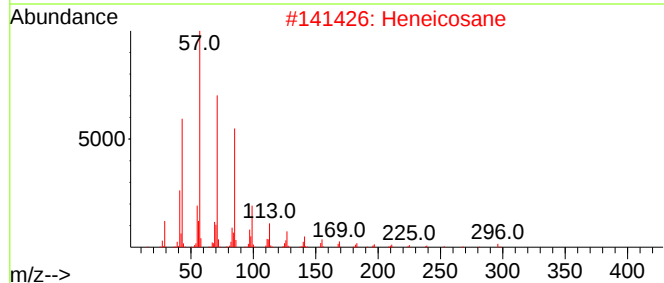
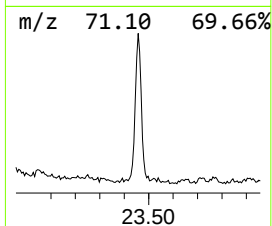
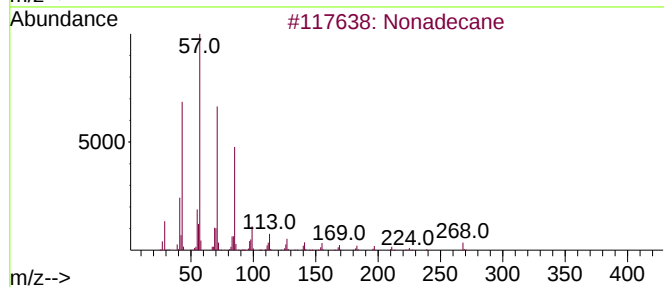
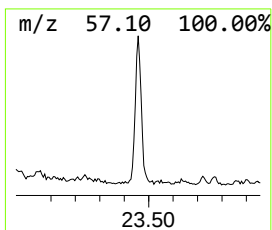
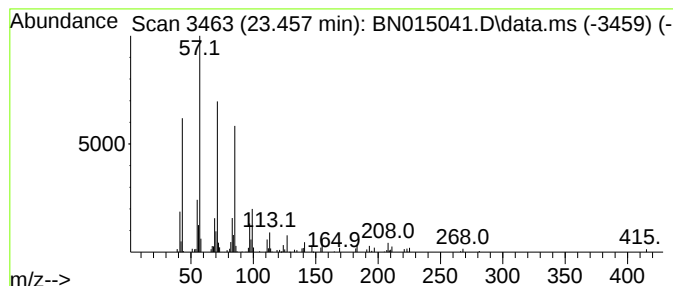
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.457	2.25 ng/ul	247337	Perylene-d12	23.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	90
2			Heneicosane	296	C21H44	000629-94-7	90
3			2-methyloctacosane	408	C29H60	1000376-72-8	87
4			Hexacosane	366	C26H54	000630-01-3	87
5			Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
Acq On : 24 Jun 2021 15:30
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Sample : M2682-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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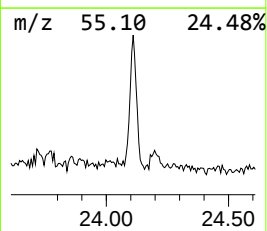
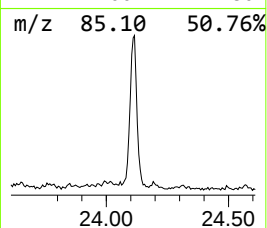
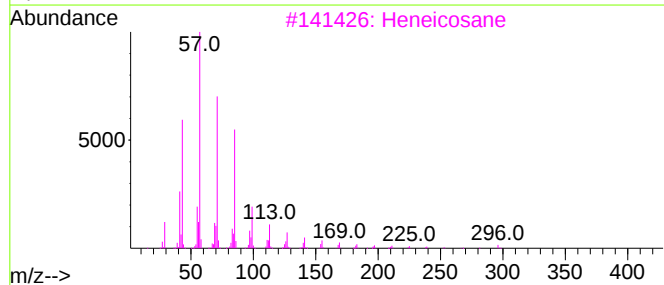
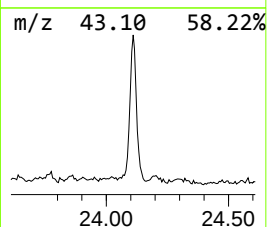
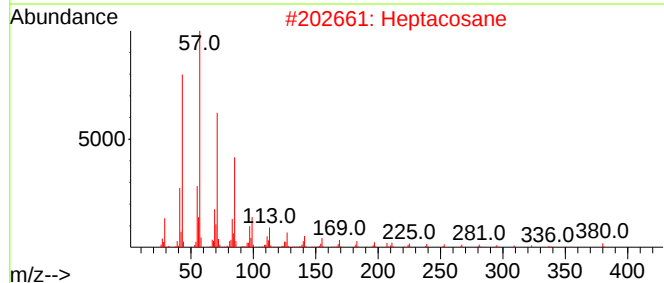
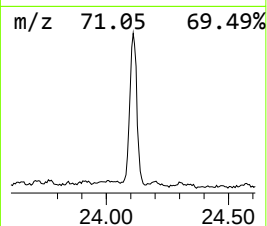
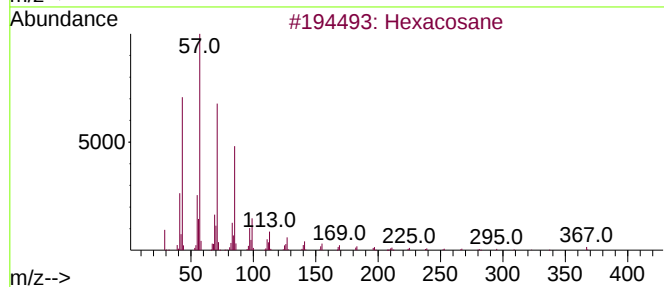
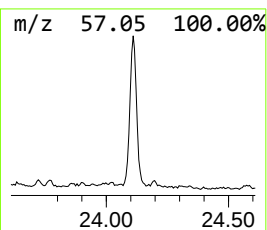
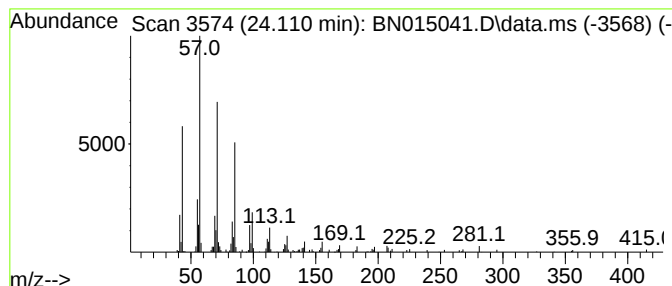
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.110	3.20 ng/ul	352226	Perylene-d12	23.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	93
2			Heptacosane	380	C27H56	000593-49-7	91
3			Heneicosane	296	C21H44	000629-94-7	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Nonadecane	268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015041.D
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Sample : M2682-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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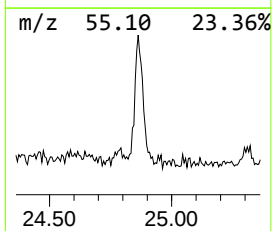
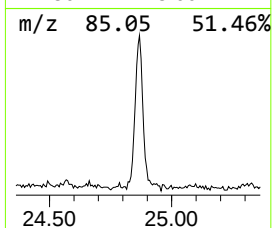
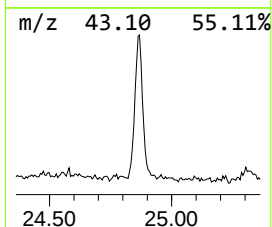
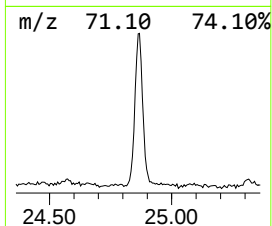
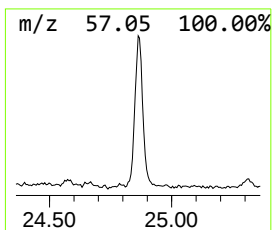
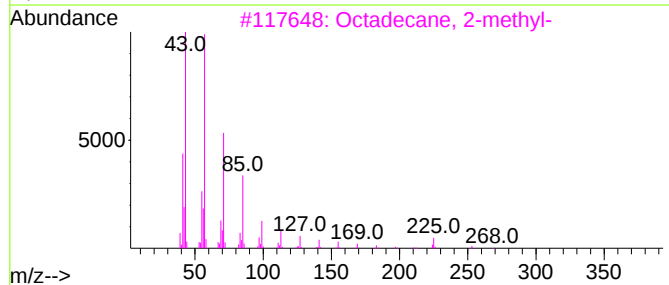
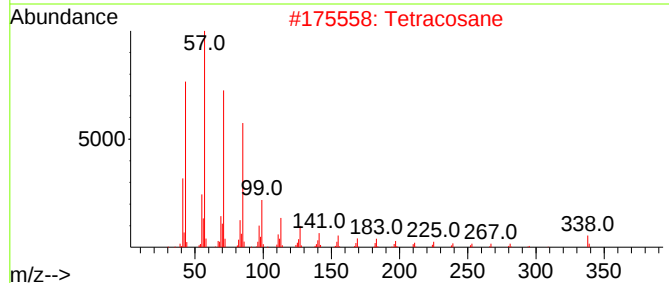
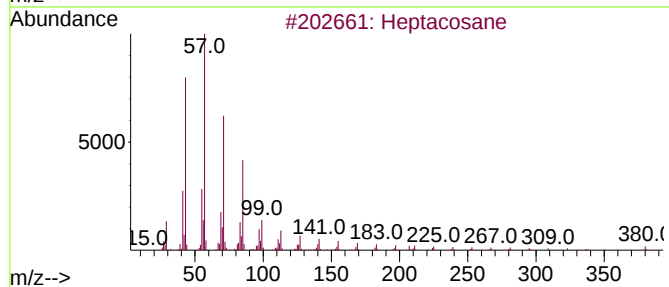
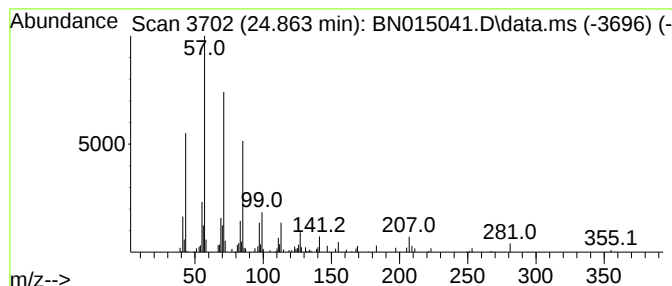
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.863	2.42 ng/ul	265509	Perylene-d12	23.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
4			2-methyloctacosane	408	C29H60	1000376-72-8	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
 Data File : BN015041.D
 Acq On : 24 Jun 2021 15:30
 Operator : CG/JU
 Sample : M2682-01
 Misc :
 ALS Vial : 5 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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unknown-01	19.086	25.0	ng/ul	2447360	4	17.098	1959770	20.0
N-Methyl-N-phen...	19.898	3.6	ng/ul	389356	5	21.298	2158270	20.0
Adipic acid, 2-...	20.522	31.6	ng/ul	3414450	5	21.298	2158270	20.0
(DEL) Alkane: S...	21.910	4.0	ng/ul	425904	5	21.298	2158270	20.0
(DEL) Alkane: S...	22.886	4.7	ng/ul	512945	6	23.539	2198260	20.0
(DEL) Alkane: S...	23.457	2.3	ng/ul	247337	6	23.539	2198260	20.0
(DEL) Alkane: S...	24.110	3.2	ng/ul	352226	6	23.539	2198260	20.0
(DEL) Alkane: S...	24.863	2.4	ng/ul	265509	6	23.539	2198260	20.0
2-Isopropenyl-4...	27.915	4.1	ng/ul	447464	6	23.539	2198260	20.0