

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
 Data File : BN015045.D
 Acq On : 24 Jun 2021 17:54
 Operator : CG/JU
 Sample : M2682-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD95

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: BN015045.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	42	rVB	29479	39527	0.39%	0.090%
2	3.681	98	101	112	rBV	194246	304759	3.02%	0.690%
3	3.787	116	119	125	rVV	25037	36708	0.36%	0.083%
4	3.910	137	140	148	rVV2	11888	19114	0.19%	0.043%
5	4.004	151	156	166	rVB	40376	60345	0.60%	0.137%
6	4.357	210	216	224	rBV5	14547	23805	0.24%	0.054%
7	4.534	242	246	255	rVB	31579	46150	0.46%	0.105%
8	5.093	336	341	347	rBV	42495	65053	0.65%	0.147%
9	5.287	368	374	380	rVB	23860	32441	0.32%	0.073%
10	5.410	391	395	405	rVB	66875	134949	1.34%	0.306%
11	5.775	452	457	464	rVB2	34401	51415	0.51%	0.116%
12	6.893	639	647	656	rBV	416604	675346	6.70%	1.530%
13	7.069	670	677	686	rBV	315287	514639	5.11%	1.166%
14	7.263	703	710	717	rVB	426241	673119	6.68%	1.525%
15	7.363	720	727	735	rBV5	8815	15729	0.16%	0.036%
16	7.728	783	789	796	rBV	651262	1008041	10.01%	2.284%
17	7.793	796	800	807	rVB	22416	32468	0.32%	0.074%
18	8.422	900	907	915	rBV	450983	736640	7.31%	1.669%
19	8.540	921	927	933	rVB	25097	40908	0.41%	0.093%
20	8.869	977	983	991	rVB	353953	572615	5.68%	1.297%
21	9.587	1098	1105	1112	rBV	250573	406270	4.03%	0.920%
22	10.116	1188	1195	1206	rVB	453463	759979	7.54%	1.722%
23	10.281	1217	1223	1230	rBV2	30003	51016	0.51%	0.116%
24	10.498	1253	1260	1267	rBV	844623	1415343	14.05%	3.207%
25	10.634	1276	1283	1296	rVB	363124	625551	6.21%	1.417%
26	11.275	1386	1392	1396	rBV3	9078	14714	0.15%	0.033%
27	11.675	1454	1460	1463	rBV2	10011	15878	0.16%	0.036%
28	12.845	1654	1659	1665	rBV2	12929	22984	0.23%	0.052%
29	13.192	1714	1718	1722	rVB2	11550	15338	0.15%	0.035%
30	13.675	1792	1800	1803	rBV5	8731	16403	0.16%	0.037%
31	13.763	1804	1815	1821	rVV	804648	1133587	11.25%	2.568%
32	13.845	1824	1829	1835	rVB6	7535	17020	0.17%	0.039%
33	14.045	1856	1863	1873	rBV	932766	1396782	13.86%	3.164%
34	14.269	1894	1901	1907	rVB3	21731	33142	0.33%	0.075%
35	14.351	1907	1915	1923	rBV	1201185	1813887	18.00%	4.109%
36	14.539	1939	1947	1954	rBV	291292	441989	4.39%	1.001%

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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

37	14.692	1968	1973	1979	rVB8	14317	26172	0.26%	0.059%
38	14.769	1979	1986	1992	rBV6	33937	67434	0.67%	0.153%
39	15.127	2043	2047	2050	rBV3	16480	21151	0.21%	0.048%
40	15.227	2060	2064	2070	rVB4	16121	21661	0.21%	0.049%

41	15.345	2074	2084	2090	rBV	1194162	1744917	17.32%	3.953%
42	15.398	2090	2093	2098	rVB4	22974	32469	0.32%	0.074%
43	15.457	2098	2103	2109	rBV	237655	329207	3.27%	0.746%
44	15.504	2109	2111	2118	rVB4	16337	24080	0.24%	0.055%
45	15.586	2121	2125	2129	rVB3	13480	17311	0.17%	0.039%

46	15.692	2136	2143	2151	rBV8	12710	35999	0.36%	0.082%
47	15.774	2151	2157	2165	rBV2	119342	194502	1.93%	0.441%
48	15.845	2166	2169	2176	rVB5	14990	22743	0.23%	0.052%
49	15.986	2188	2193	2196	rBV6	6459	14638	0.15%	0.033%
50	16.022	2196	2199	2205	rVB7	12146	21463	0.21%	0.049%

51	16.086	2205	2210	2215	rBV4	34847	48665	0.48%	0.110%
52	16.139	2215	2219	2223	rVV5	17180	27120	0.27%	0.061%
53	16.310	2244	2248	2254	rBV7	12934	18356	0.18%	0.042%
54	16.392	2257	2262	2269	rBV3	70638	112913	1.12%	0.256%
55	16.651	2301	2306	2308	rBV4	15666	22786	0.23%	0.052%

56	16.833	2332	2337	2342	rBV7	21961	45451	0.45%	0.103%
57	16.880	2342	2345	2349	rVB2	32700	41743	0.41%	0.095%
58	17.098	2376	2382	2387	rBV	1415257	2096874	20.81%	4.751%
59	17.139	2387	2389	2393	rVV	68512	88506	0.88%	0.201%
60	17.198	2393	2399	2409	rVV	1312646	1912828	18.99%	4.334%

61	17.298	2411	2416	2421	rVB8	23820	41434	0.41%	0.094%
62	17.621	2467	2471	2475	rVV2	42280	57812	0.57%	0.131%
63	17.786	2495	2499	2505	rBV7	29816	52032	0.52%	0.118%
64	17.892	2514	2517	2521	rVB6	12570	15120	0.15%	0.034%
65	17.980	2528	2532	2533	rBV	19723	25881	0.26%	0.059%

66	18.004	2533	2536	2544	rVV2	118926	209775	2.08%	0.475%
67	18.074	2546	2548	2552	rVB	12598	16304	0.16%	0.037%
68	18.168	2561	2564	2568	rVV5	19223	24750	0.25%	0.056%
69	18.210	2568	2571	2574	rVB3	16763	21383	0.21%	0.048%
70	18.316	2584	2589	2596	rBV	49772	74968	0.74%	0.170%

71	18.427	2603	2608	2613	rBV	407568	524391	5.20%	1.188%
72	18.721	2656	2658	2662	rVB5	19968	23874	0.24%	0.054%
73	18.780	2665	2668	2671	rBV3	14736	17080	0.17%	0.039%
74	18.910	2684	2690	2693	rBV7	30510	67566	0.67%	0.153%
75	18.951	2693	2697	2702	rVV2	72835	110159	1.09%	0.250%

76	19.163	2729	2733	2740	rVB	137417	191714	1.90%	0.434%
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77	19.292	2752	2755	2762	rVB6	47720	70026	0.70%	0.159%
78	19.498	2784	2790	2794	rBV	1556865	2269739	22.53%	5.142%
79	19.857	2847	2851	2856	rBV6	16706	38369	0.38%	0.087%
80	20.068	2884	2887	2890	rBV	81362	78710	0.78%	0.178%
81	20.521	2959	2964	2970	rBV	8493121	10074930	100.00%	22.825%
82	20.862	3019	3022	3025	rBV2	35598	42956	0.43%	0.097%
83	21.027	3045	3050	3055	rBV3	194638	324663	3.22%	0.736%
84	21.227	3080	3084	3089	rVB	172505	194054	1.93%	0.440%
85	21.298	3089	3096	3100	rBV	1684506	2345465	23.28%	5.314%
86	21.421	3114	3117	3120	rBV5	44595	61818	0.61%	0.140%
87	21.904	3196	3199	3204	rBV5	90364	172047	1.71%	0.390%
88	22.045	3220	3223	3238	rVB7	68476	180523	1.79%	0.409%
89	22.162	3239	3243	3247	rBV	114924	151195	1.50%	0.343%
90	22.345	3271	3274	3276	rBV3	93981	117982	1.17%	0.267%
91	22.798	3347	3351	3358	rVB2	217822	314227	3.12%	0.712%
92	22.880	3361	3365	3373	rVB2	160163	274173	2.72%	0.621%
93	23.398	3447	3453	3459	rBV	968047	1802108	17.89%	4.083%
94	23.456	3459	3463	3470	rVB3	182012	314203	3.12%	0.712%
95	23.545	3471	3478	3491	rVB2	1277010	2421333	24.03%	5.486%
96	23.721	3504	3508	3514	rVB3	72317	119933	1.19%	0.272%
97	23.886	3532	3536	3543	rVB5	104495	223688	2.22%	0.507%
98	24.109	3568	3574	3579	rBV	163027	312098	3.10%	0.707%
99	24.862	3696	3702	3709	rVB2	141341	317762	3.15%	0.720%
100	25.750	3848	3853	3861	rVB3	79033	190676	1.89%	0.432%

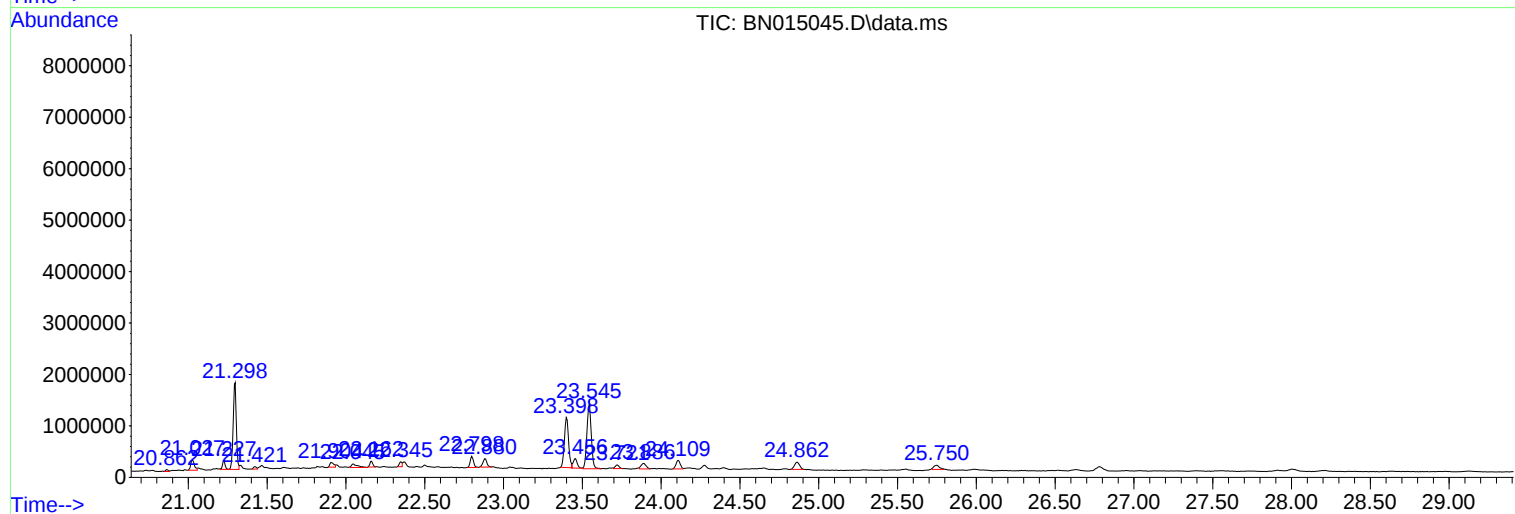
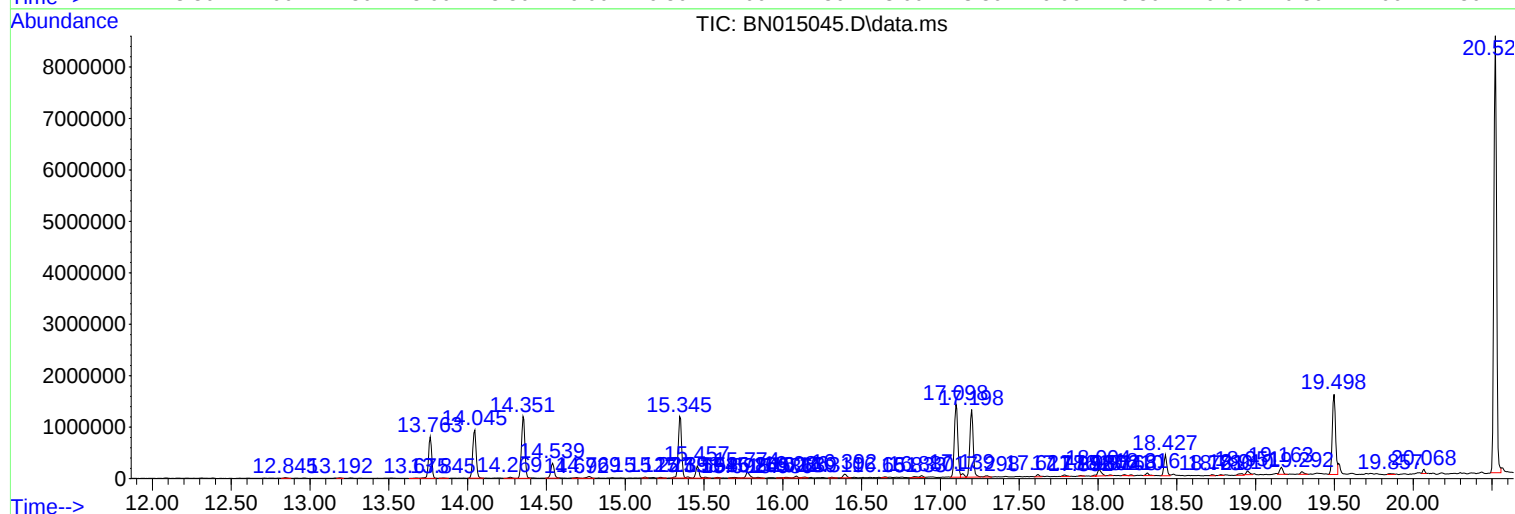
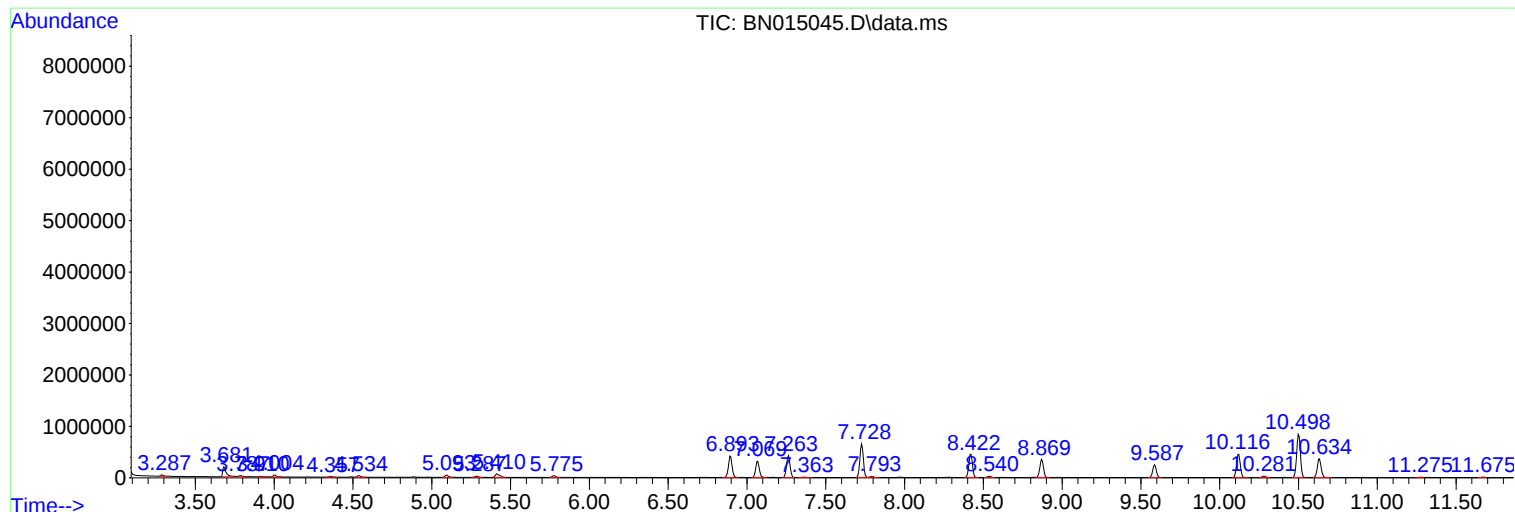
Sum of corrected areas: 44139564

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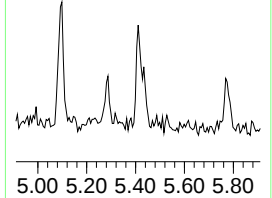
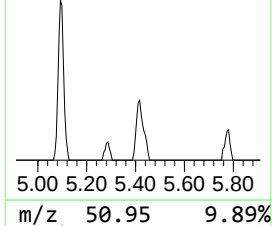
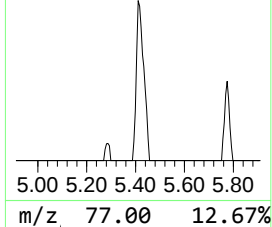
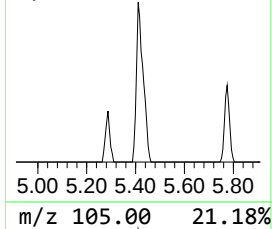
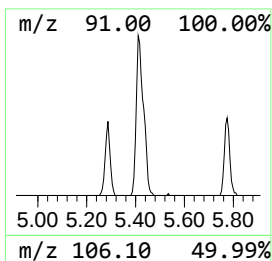
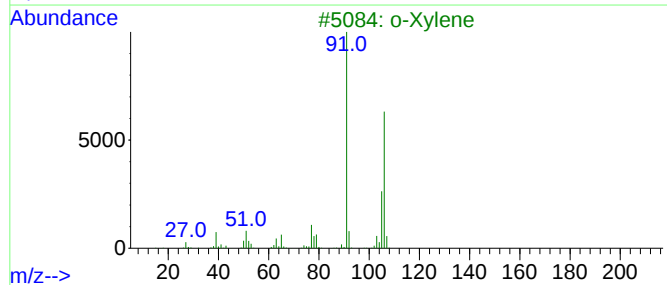
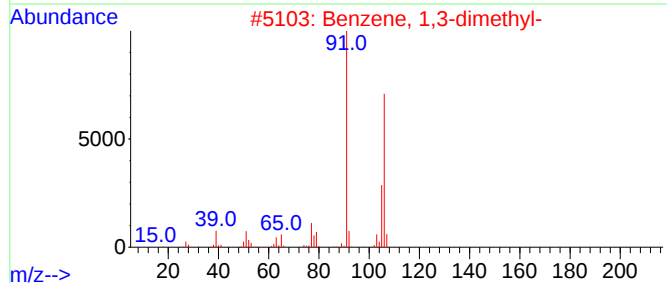
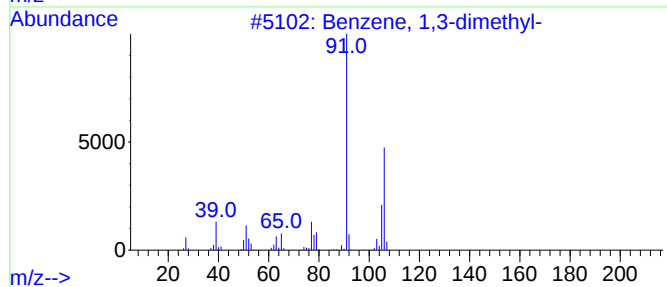
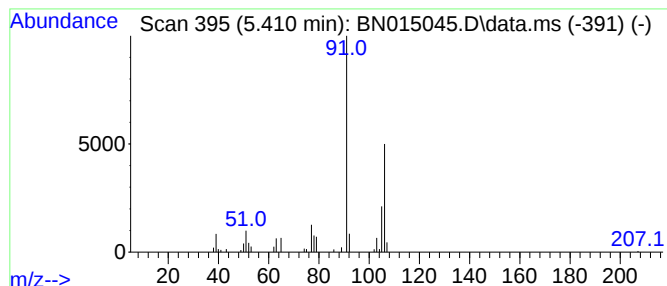
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.410	2.68 ng/ul	134949	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			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
3			o-Xylene	106	C8H10	000095-47-6	93
4			p-Xylene	106	C8H10	000106-42-3	91
5			o-Xylene	106	C8H10	000095-47-6	91



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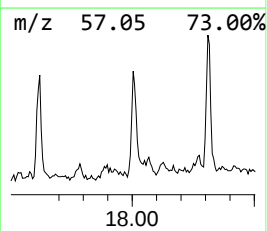
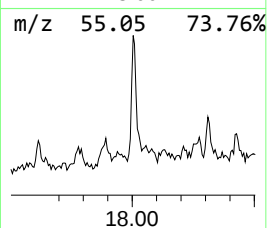
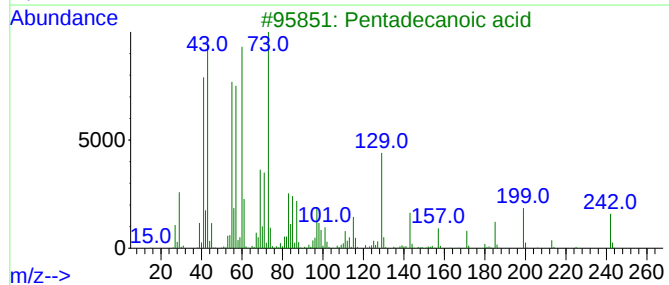
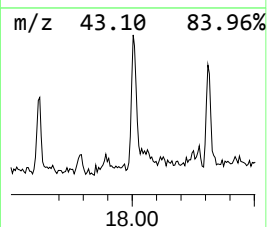
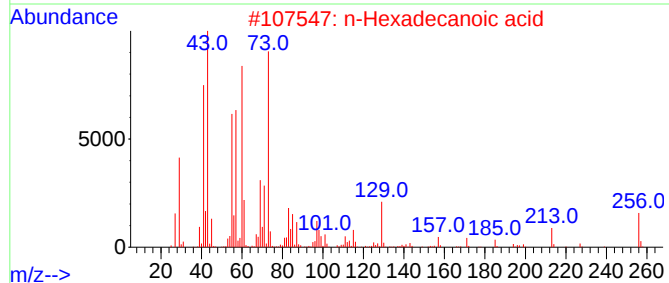
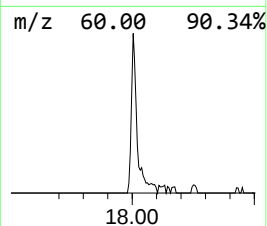
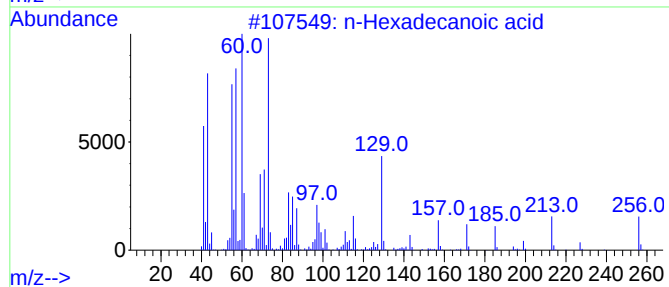
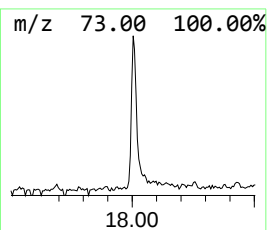
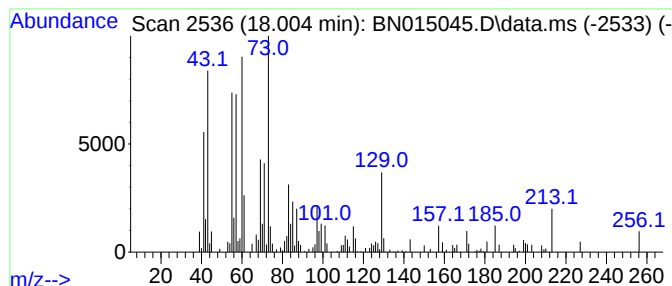
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.004	2.00 ng/ul	209775	Phenanthrene-d10	17.098

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	97
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			Dodecanoic acid	200	C12H24O2	000143-07-7	70



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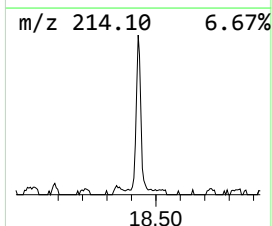
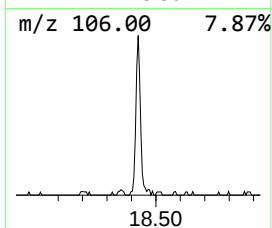
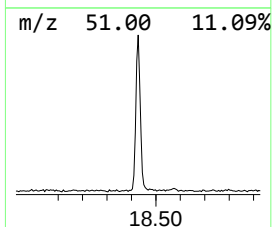
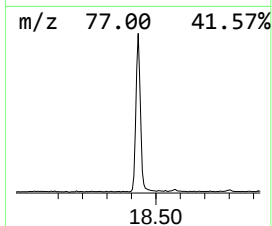
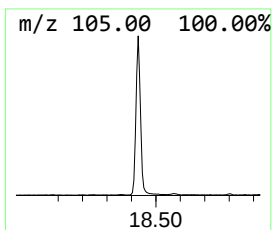
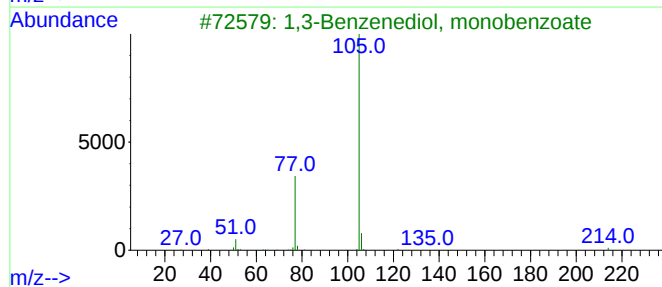
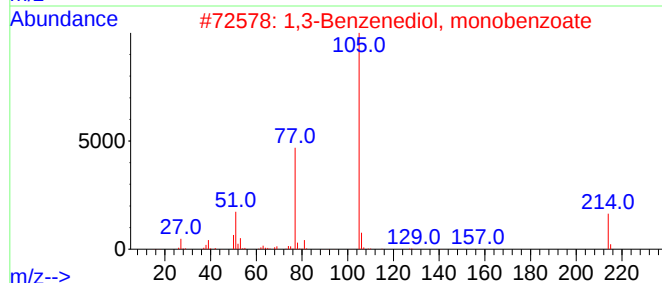
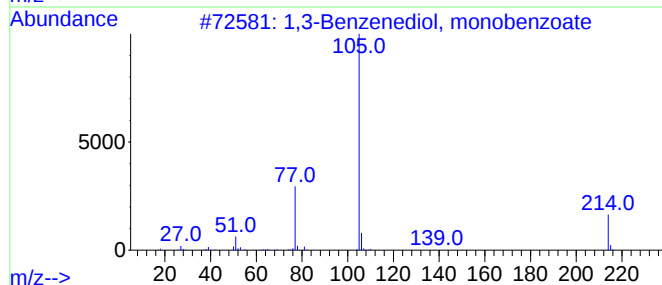
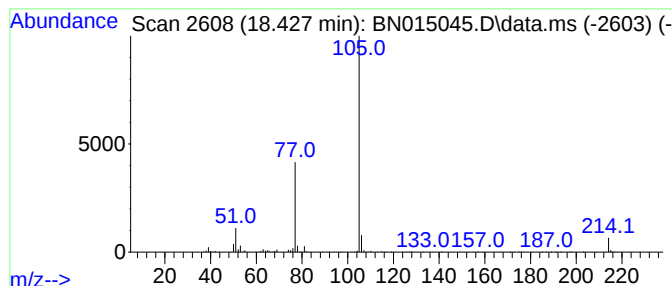
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,3-Benzenediol, monobenzoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.427	5.00 ng/ul	524391	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	91		
2	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90		
3	1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90		
4	Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	80		
5	1-Benzoyloxy-2-formyloxybenzene	242	C14H10O4	079792-93-1	78		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
Acq On : 24 Jun 2021 17:54
Operator : CG/JU
Sample : M2682-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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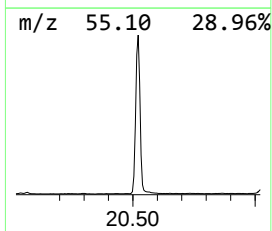
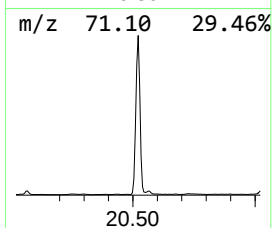
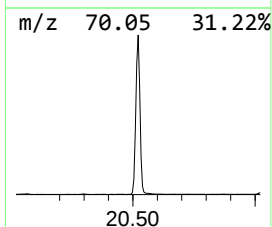
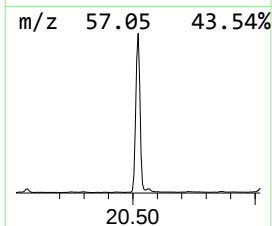
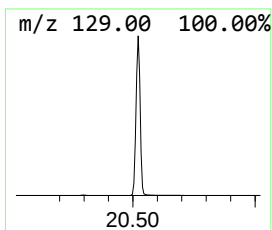
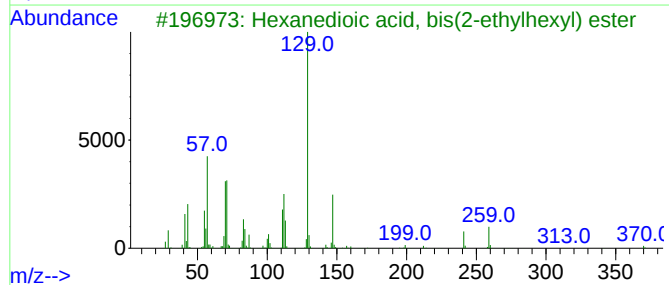
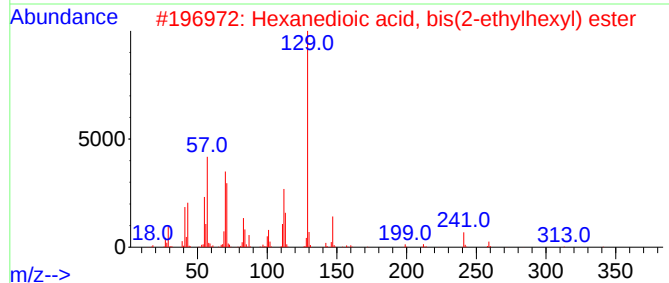
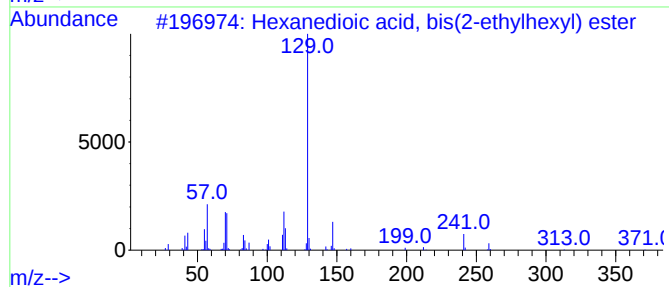
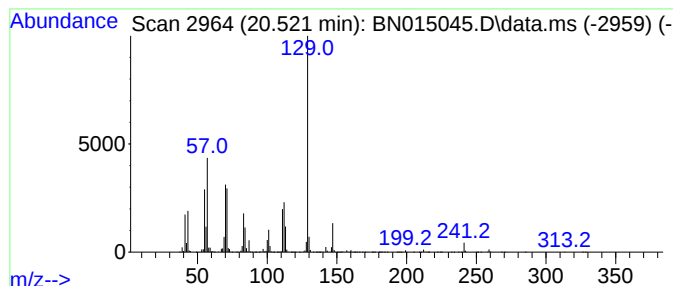
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.521	85.91 ng/ul	10074900	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4			Diisooctyl adipate	370	C22H42O4	001330-86-5	83
5			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
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Sample : M2682-08
Misc :
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ClientSampleId :
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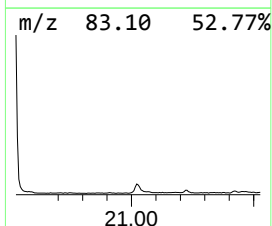
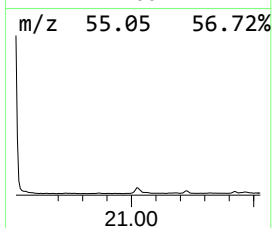
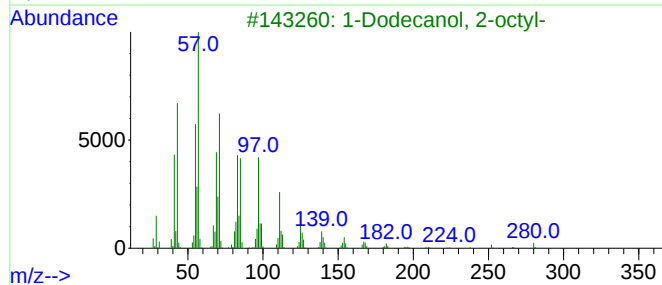
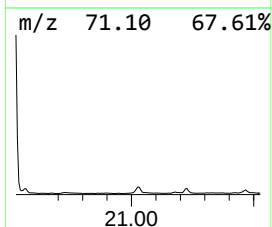
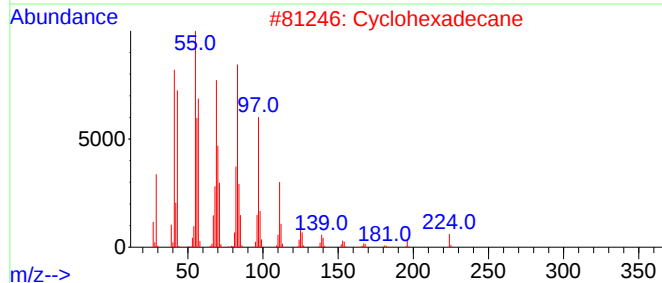
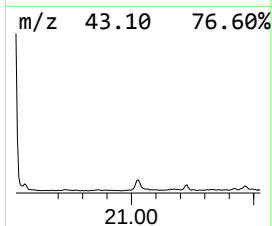
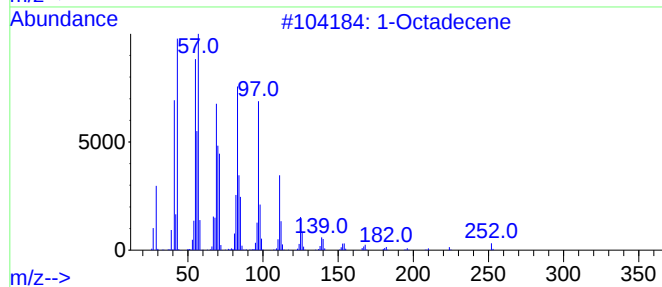
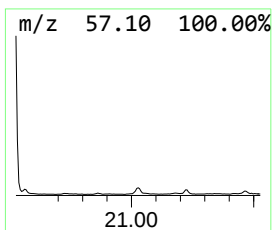
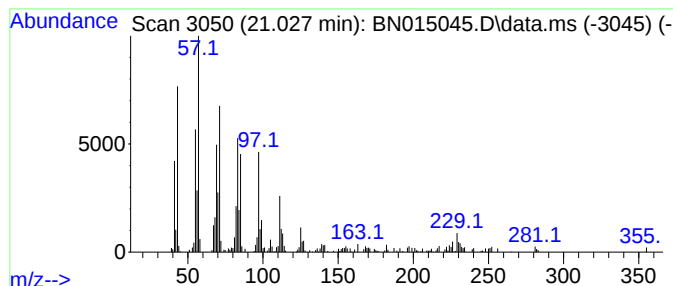
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	2.77 ng/ul	324663	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecene	252	C18H36	000112-88-9	98
2			Cyclohexadecane	224	C16H32	000295-65-8	95
3			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	91
4			1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	91
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
Acq On : 24 Jun 2021 17:54
Operator : CG/JU
Sample : M2682-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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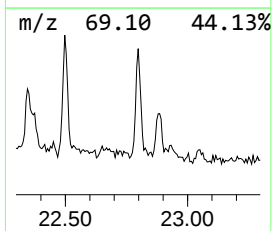
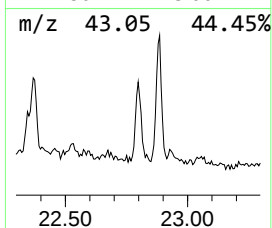
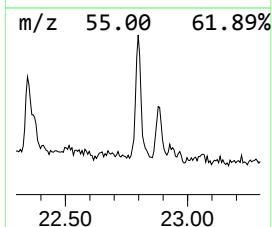
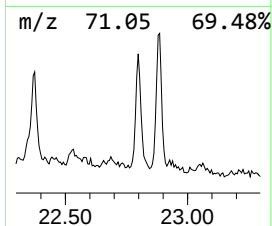
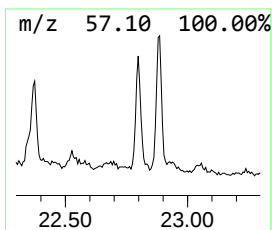
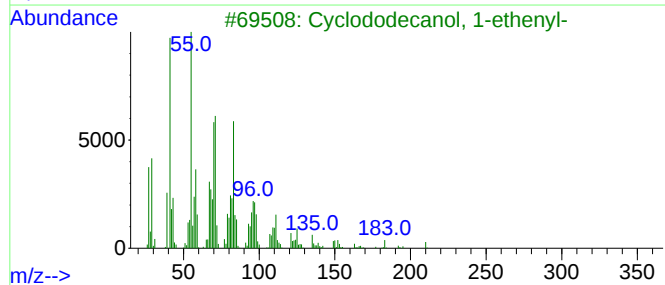
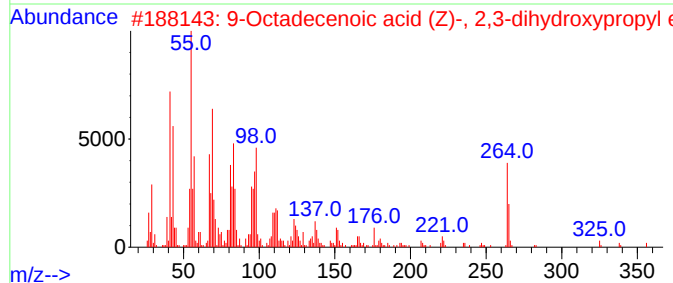
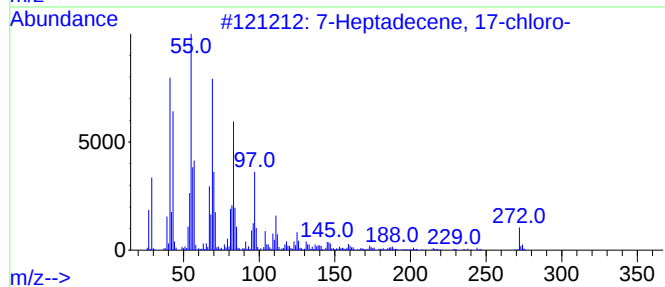
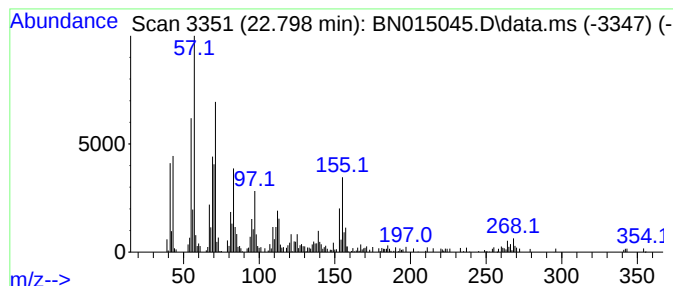
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 7-Heptadecene, 17-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.798	2.60 ng/ul	314227	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Heptadecene, 17-chloro-	272	C17H33Cl	056554-79-1	55
2			9-Octadecenoic acid (Z)-, 2,3-di...	356	C21H40O4	000111-03-5	45
3			Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	43
4			Oxalic acid, decyl 2-ethylhexyl ...	342	C20H38O4	1000309-39-3	43
5			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
Acq On : 24 Jun 2021 17:54
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Sample : M2682-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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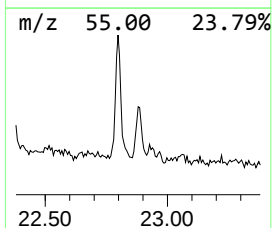
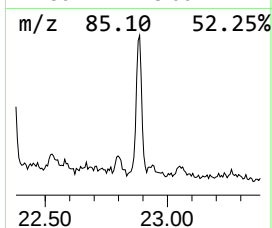
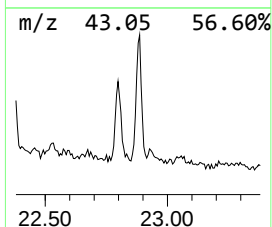
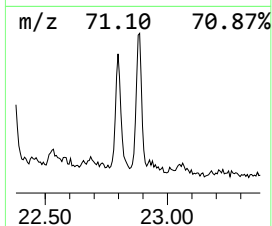
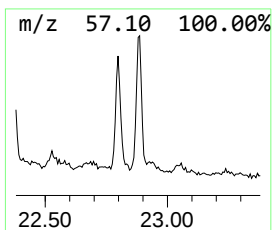
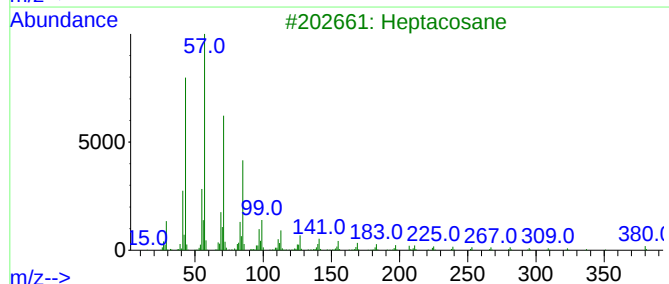
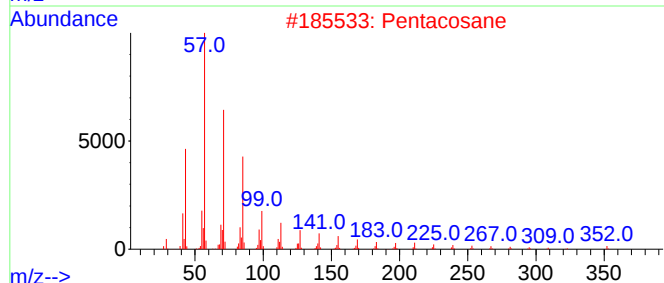
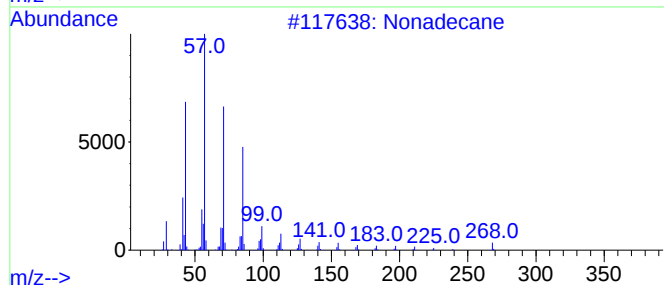
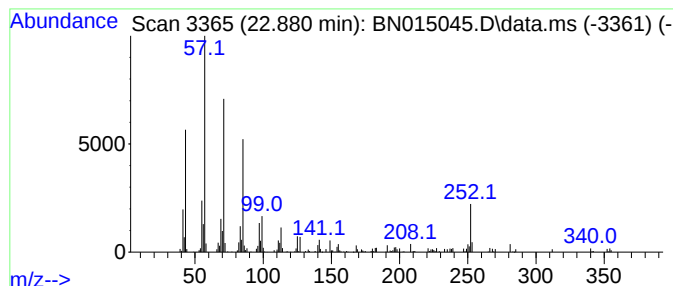
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.880	2.26 ng/ul	274173	Perylene-d12	23.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	93
2		Pentacosane	352	C25H52	000629-99-2	90
3		Heptacosane	380	C27H56	000593-49-7	81
4		Nonadecane	268	C19H40	000629-92-5	76
5		Heneicosane	296	C21H44	000629-94-7	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
Acq On : 24 Jun 2021 17:54
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Sample : M2682-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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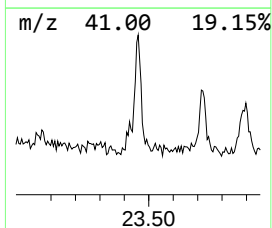
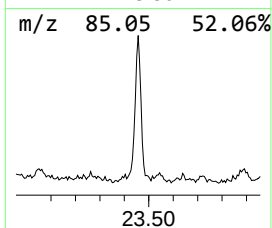
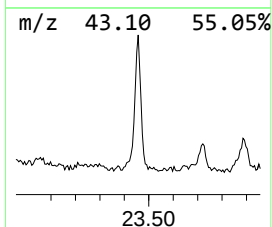
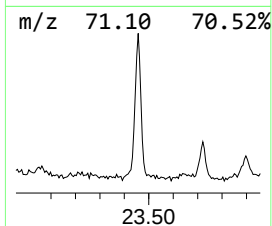
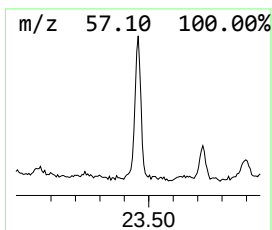
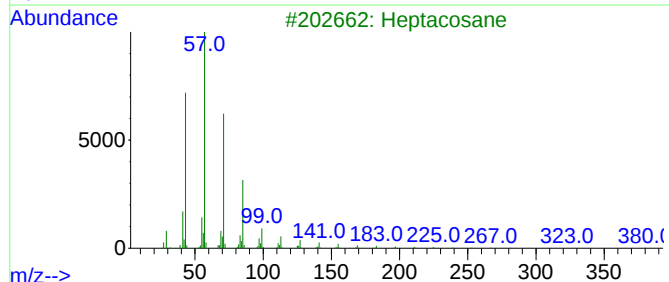
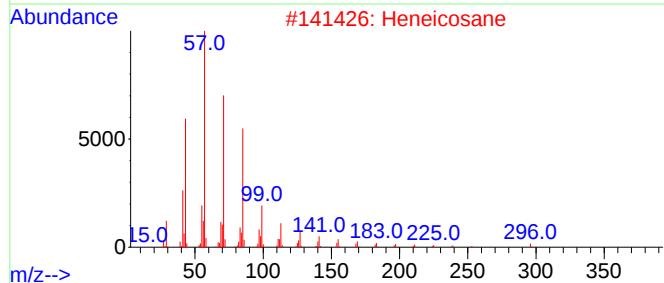
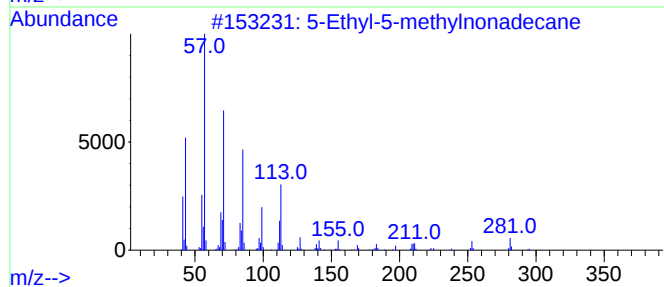
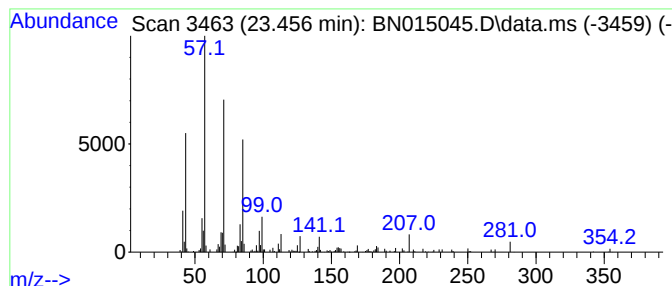
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.456	2.60 ng/ul	314203	Perylene-d12	23.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	86
2		Heneicosane	296	C21H44	000629-94-7	83
3		Heptacosane	380	C27H56	000593-49-7	80
4		Heptadecane	240	C17H36	000629-78-7	80
5		triacontane	422	C30H62	000638-68-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
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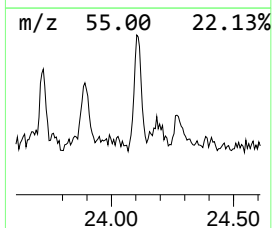
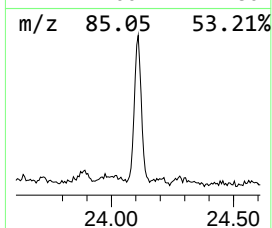
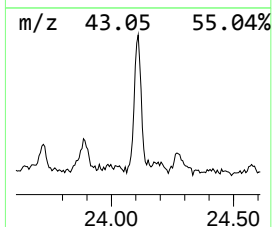
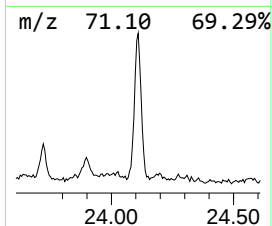
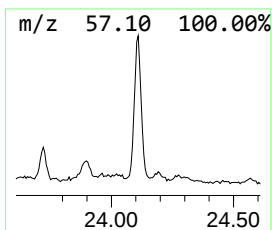
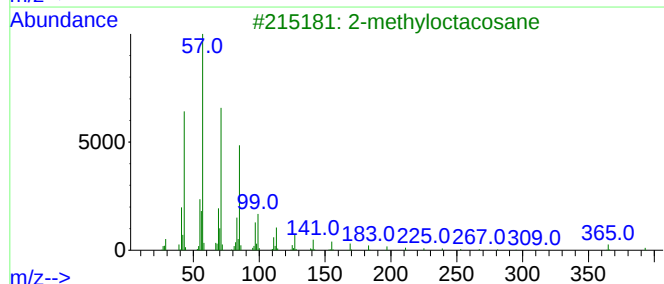
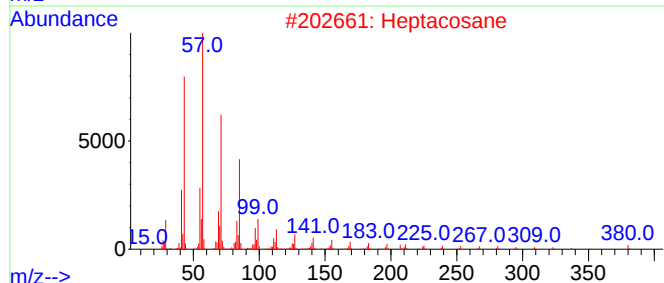
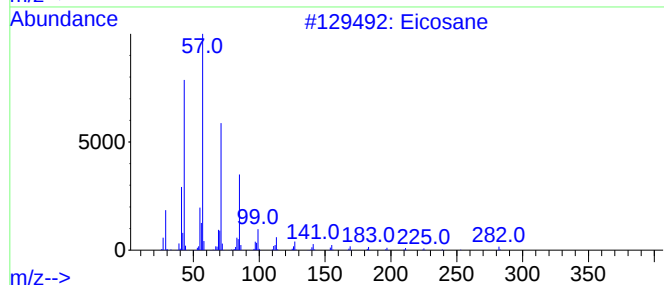
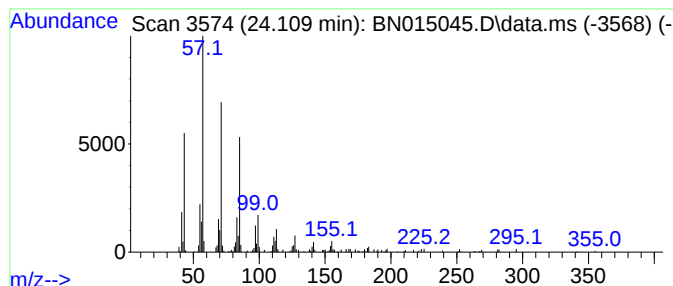
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Quant Title : SVOA CALIBRATION

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.
24.109	2.58 ng/ul	312098	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Heptacosane	380	C27H56	000593-49-7	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	90
4			Tetracosane	338	C24H50	000646-31-1	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015045.D
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Sample : M2682-08
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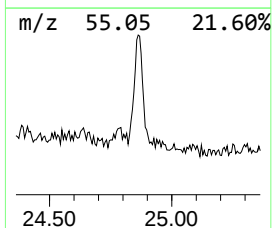
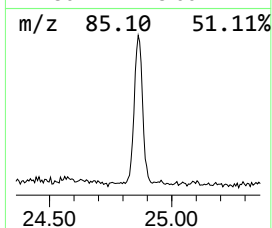
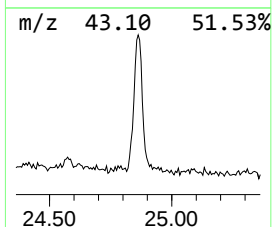
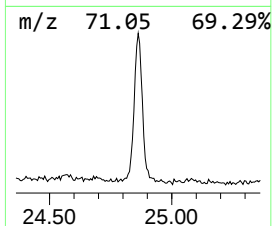
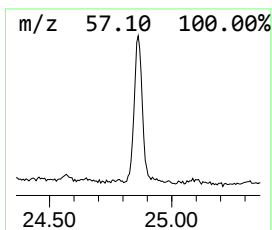
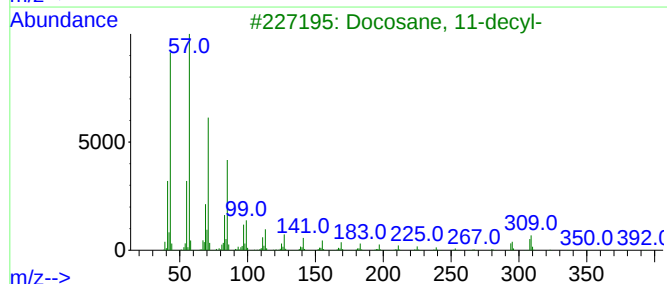
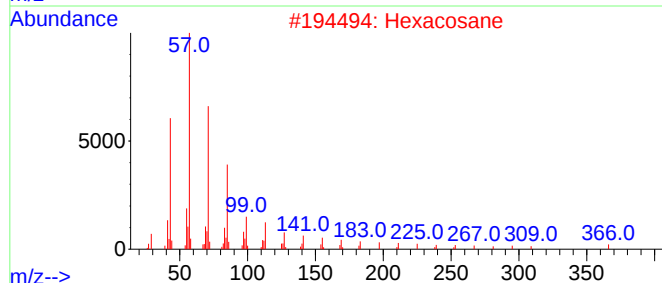
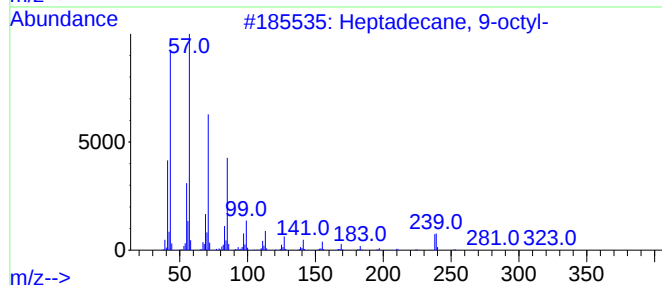
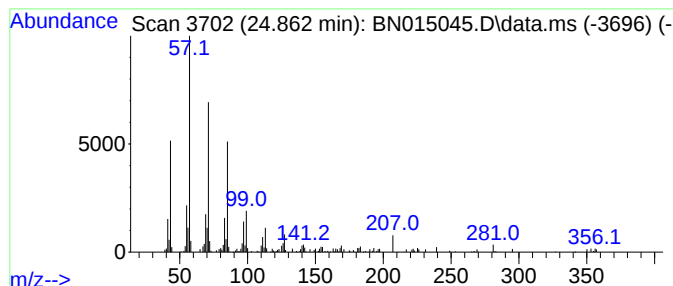
Instrument :
BNA_N
ClientSampleId :
BGD95

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.862	2.62 ng/ul	317762	Perylene-d12	23.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Hexacosane	366	C26H54	000630-01-3	86
3	Docosane, 11-decyl-	451	C32H66	055401-55-3	86
4	Pentacosane	352	C25H52	000629-99-2	86
5	Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
 Data File : BN015045.D
 Acq On : 24 Jun 2021 17:54
 Operator : CG/JU
 Sample : M2682-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD95

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
Benzene, 1,3-di...	5.410	2.7	ng/ul	134949	1	7.728	1008040	20.0
n-Hexadecanoic ...	18.004	2.0	ng/ul	209775	4	17.098	2096870	20.0
1,3-Benzenediol...	18.427	5.0	ng/ul	524391	4	17.098	2096870	20.0
Hexanedioic aci...	20.521	85.9	ng/ul	10074900	5	21.298	2345470	20.0
(DEL) Alkane: S...	21.027	2.8	ng/ul	324663	5	21.298	2345470	20.0
7-Heptadecene, ...	22.798	2.6	ng/ul	314227	6	23.545	2421330	20.0
(DEL) Alkane: S...	22.880	2.3	ng/ul	274173	6	23.545	2421330	20.0
(DEL) Alkane: S...	23.456	2.6	ng/ul	314203	6	23.545	2421330	20.0
(DEL) Alkane: S...	24.109	2.6	ng/ul	312098	6	23.545	2421330	20.0
(DEL) Alkane: S...	24.862	2.6	ng/ul	317762	6	23.545	2421330	20.0