

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
 Data File : BG051872.D
 Acq On : 4 Jan 2022 16:14
 Operator : CG/JU
 Sample : M5215-01MEDL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLD3MEDL

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_G\Methods\SFAM-EPA-BG121421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051872.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.726	377	381	388	rVV	46969	72299	7.94%	1.178%
2	5.862	398	404	415	rVB	109448	212924	23.38%	3.471%
3	6.243	463	469	476	rBV2	30983	49898	5.48%	0.813%
4	6.502	508	513	520	rBV6	2876	6057	0.67%	0.099%
5	6.790	558	562	567	rBV2	4282	6517	0.72%	0.106%
6	6.878	573	577	580	rVV4	3989	6154	0.68%	0.100%
7	7.224	634	636	643	rVV4	6379	10941	1.20%	0.178%
8	7.289	643	647	650	rVV3	4539	6796	0.75%	0.111%
9	7.336	650	655	660	rVV3	9688	14523	1.59%	0.237%
10	7.395	660	665	673	rVV3	14071	25142	2.76%	0.410%
11	7.477	673	679	683	rVB5	5734	9745	1.07%	0.159%
12	7.565	689	694	701	rBV3	5348	10498	1.15%	0.171%
13	7.641	701	707	711	rVV3	4764	6353	0.70%	0.104%
14	7.788	726	732	736	rVV3	4944	9767	1.07%	0.159%
15	7.871	740	746	749	rVV2	18302	30259	3.32%	0.493%
16	7.900	749	751	757	rVB2	16184	21723	2.39%	0.354%
17	8.258	803	812	819	rBV	108083	195975	21.52%	3.194%
18	8.382	828	833	839	rVB3	4636	7709	0.85%	0.126%
19	8.916	917	924	928	rVV5	4275	10049	1.10%	0.164%
20	8.963	928	932	938	rVB3	4038	6604	0.73%	0.108%
21	9.427	1006	1011	1018	rVV3	4804	9136	1.00%	0.149%
22	9.492	1018	1022	1027	rVB3	13534	20635	2.27%	0.336%
23	10.162	1130	1136	1141	rBV4	4529	7794	0.86%	0.127%
24	10.708	1222	1229	1234	rBV2	4652	7869	0.86%	0.128%
25	11.054	1281	1288	1289	rBV3	7472	10380	1.14%	0.169%
26	11.096	1289	1295	1301	rVV	154675	279823	30.73%	4.561%
27	11.143	1301	1303	1311	rVB2	8418	12594	1.38%	0.205%
28	12.100	1460	1466	1470	rVB3	3089	6010	0.66%	0.098%
29	12.488	1527	1532	1537	rVB2	11830	17446	1.92%	0.284%
30	12.740	1570	1575	1580	rVB3	5276	8337	0.92%	0.136%
31	13.428	1687	1692	1696	rVB2	5821	8186	0.90%	0.133%
32	13.710	1733	1740	1750	rBV	32416	50714	5.57%	0.827%
33	14.221	1822	1827	1831	rBV6	5432	9100	1.00%	0.148%
34	14.280	1833	1837	1842	rVB3	17506	26633	2.92%	0.434%
35	14.362	1848	1851	1855	rVV3	10794	13648	1.50%	0.222%
36	14.591	1884	1890	1896	rBV2	13781	21468	2.36%	0.350%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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37	14.767	1916	1920	1927	rVB	40845	54280	5.96%	0.885%
38	14.896	1935	1942	1948	rVB2	277418	440511	48.37%	7.180%
39	15.096	1969	1976	1977	rBV5	3178	6327	0.69%	0.103%
40	15.278	2002	2007	2011	rBV7	7340	12332	1.35%	0.201%
41	15.384	2020	2025	2032	rVB7	6644	15714	1.73%	0.256%
42	15.507	2042	2046	2051	rBV6	3602	6803	0.75%	0.111%
43	15.719	2078	2082	2086	rVB	48609	60252	6.62%	0.982%
44	15.824	2096	2100	2104	rBV3	4471	6294	0.69%	0.103%
45	15.883	2105	2110	2115	rBV	24428	30001	3.29%	0.489%
46	16.065	2138	2141	2144	rVB2	4676	5740	0.63%	0.094%
47	16.124	2144	2151	2154	rBV2	18077	33444	3.67%	0.545%
48	16.224	2163	2168	2169	rBV5	6833	9730	1.07%	0.159%
49	16.283	2173	2178	2183	rVV7	5266	9783	1.07%	0.159%
50	16.341	2183	2188	2191	rVV5	7186	10653	1.17%	0.174%
51	16.377	2191	2194	2197	rBV3	5686	6488	0.71%	0.106%
52	16.576	2224	2228	2231	rBV	55536	74763	8.21%	1.219%
53	16.606	2231	2233	2238	rVB3	34377	42071	4.62%	0.686%
54	16.706	2246	2250	2252	rBV2	13026	16514	1.81%	0.269%
55	16.852	2272	2275	2279	rVB5	6297	6769	0.74%	0.110%
56	16.894	2279	2282	2285	rVV4	4509	6408	0.70%	0.104%
57	16.917	2285	2286	2290	rVV4	5020	5655	0.62%	0.092%
58	16.988	2293	2298	2299	rBV4	8651	11289	1.24%	0.184%
59	17.005	2299	2301	2304	rVV4	8602	10088	1.11%	0.164%
60	17.035	2304	2306	2310	rVB5	4831	5696	0.63%	0.093%
61	17.087	2312	2315	2321	rVB8	5319	7527	0.83%	0.123%
62	17.369	2359	2363	2368	rBV	41490	53123	5.83%	0.866%
63	17.428	2368	2373	2377	rBV2	22013	28837	3.17%	0.470%
64	17.522	2385	2389	2393	rBV2	10730	14677	1.61%	0.239%
65	17.640	2403	2409	2415	rBV	366140	571481	62.76%	9.315%
66	17.740	2421	2426	2432	rVB2	24343	40314	4.43%	0.657%
67	17.910	2451	2455	2457	rBV4	6379	9258	1.02%	0.151%
68	18.033	2474	2476	2480	rVB5	6959	7712	0.85%	0.126%
69	18.116	2486	2490	2494	rVB	45939	52240	5.74%	0.852%
70	18.280	2516	2518	2522	rVB5	8251	11216	1.23%	0.183%
71	18.374	2530	2534	2539	rBV8	5578	10820	1.19%	0.176%
72	18.468	2547	2550	2553	rBV5	4289	6560	0.72%	0.107%
73	18.503	2553	2556	2562	rBV5	24176	39213	4.31%	0.639%
74	18.556	2562	2565	2567	rBV4	6663	8347	0.92%	0.136%
75	18.797	2603	2606	2611	rVB2	33537	39170	4.30%	0.638%
76	19.073	2649	2653	2657	rVB6	12157	17772	1.95%	0.290%

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Integration Parameters: LSCINT.P

Integrator: RTE

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 Title : SVOA CALIBRATION

77	19.196	2670	2674	2676	rBV4	6640	6788	0.75%	0.111%
78	19.255	2682	2684	2689	rVB5	5552	7059	0.78%	0.115%
79	19.420	2707	2712	2716	rBV3	27442	38759	4.26%	0.632%
80	19.608	2740	2744	2748	rVB6	12643	15832	1.74%	0.258%
81	19.672	2751	2755	2758	rVB4	18517	23394	2.57%	0.381%
82	19.890	2790	2792	2795	rVB3	10433	8377	0.92%	0.137%
83	19.995	2806	2810	2811	rBV2	25078	26656	2.93%	0.434%
84	20.013	2811	2813	2818	rVB2	26906	35141	3.86%	0.573%
85	20.172	2837	2840	2841	rBV3	7094	8311	0.91%	0.135%
86	20.289	2857	2860	2863	rBV3	13159	14584	1.60%	0.238%
87	20.524	2896	2900	2905	rBV	27302	33108	3.64%	0.540%
88	20.918	2961	2967	2972	rVB	372818	502348	55.16%	8.188%
89	21.029	2982	2986	2989	rBV2	24176	27941	3.07%	0.455%
90	21.558	3073	3076	3080	rVB3	28276	37217	4.09%	0.607%
91	21.822	3115	3121	3128	rVB	631496	910642	100.00%	14.843%
92	21.957	3137	3144	3150	rVB	334081	578853	63.57%	9.435%
93	22.145	3173	3176	3181	rVB3	28509	38035	4.18%	0.620%
94	22.803	3283	3288	3296	rVB4	22029	40141	4.41%	0.654%
95	23.126	3339	3343	3352	rVB2	17058	35739	3.92%	0.583%
96	23.555	3412	3416	3422	rVB4	19096	36400	4.00%	0.593%
97	24.430	3562	3565	3570	rVB5	17501	27989	3.07%	0.456%
98	25.194	3689	3695	3700	rBV7	11918	29276	3.21%	0.477%
99	25.435	3727	3736	3749	rVB2	196319	609848	66.97%	9.941%
100	30.622	4618	4619	4622	rBV3	6130	6951	0.76%	0.113%

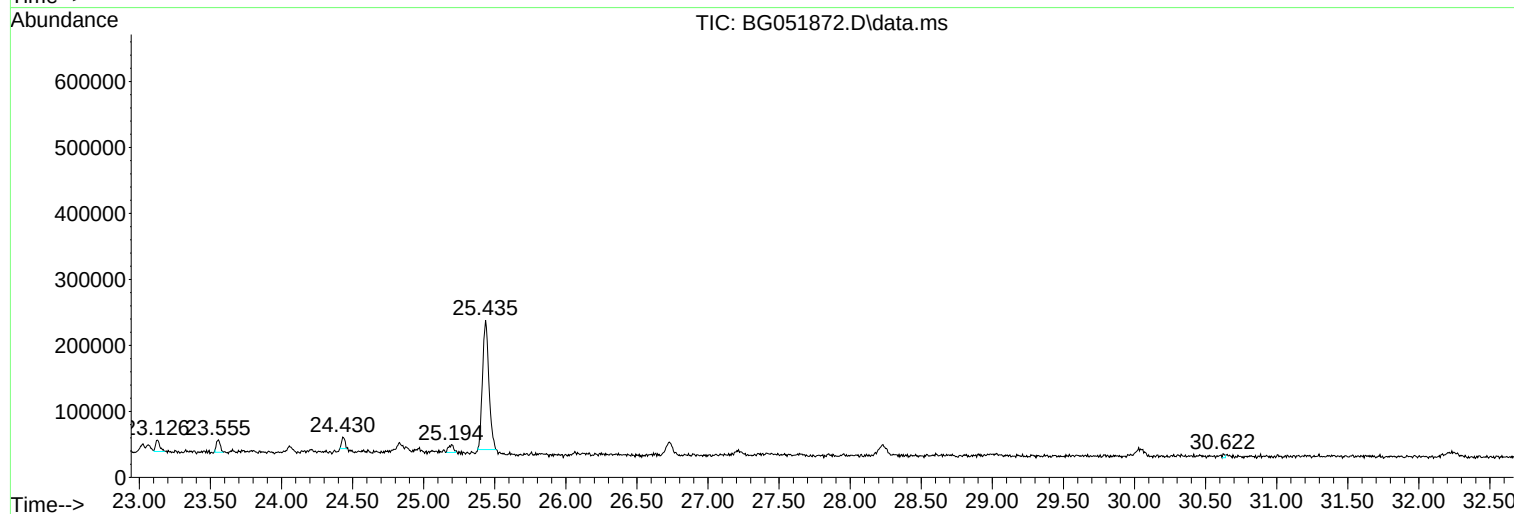
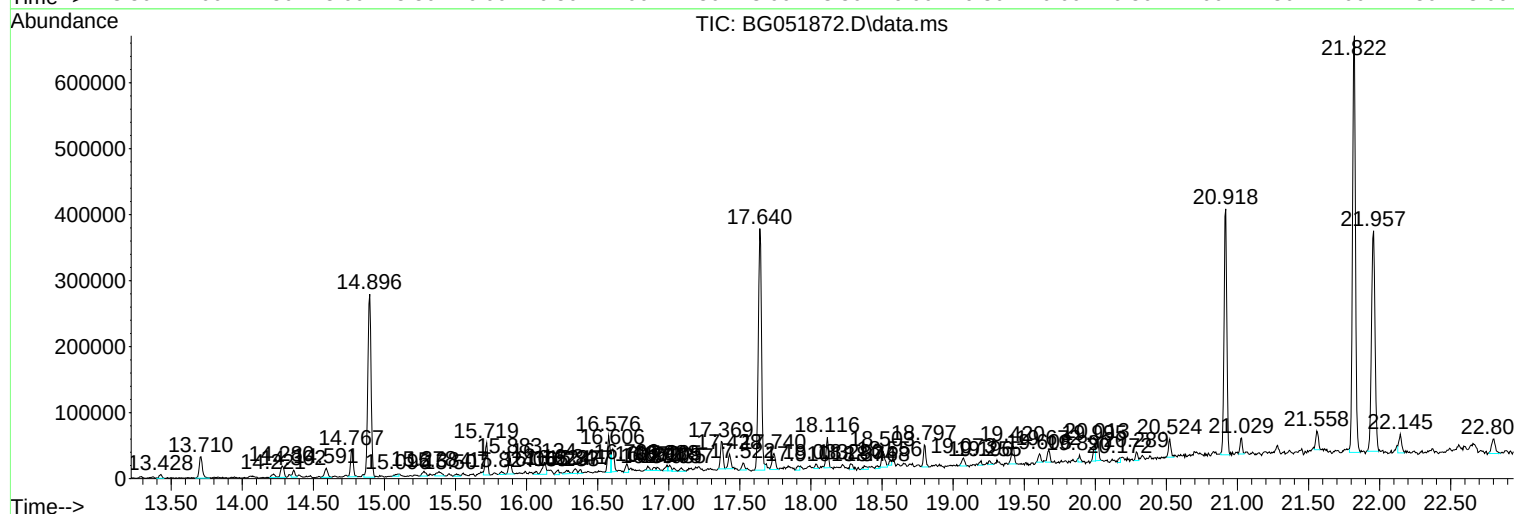
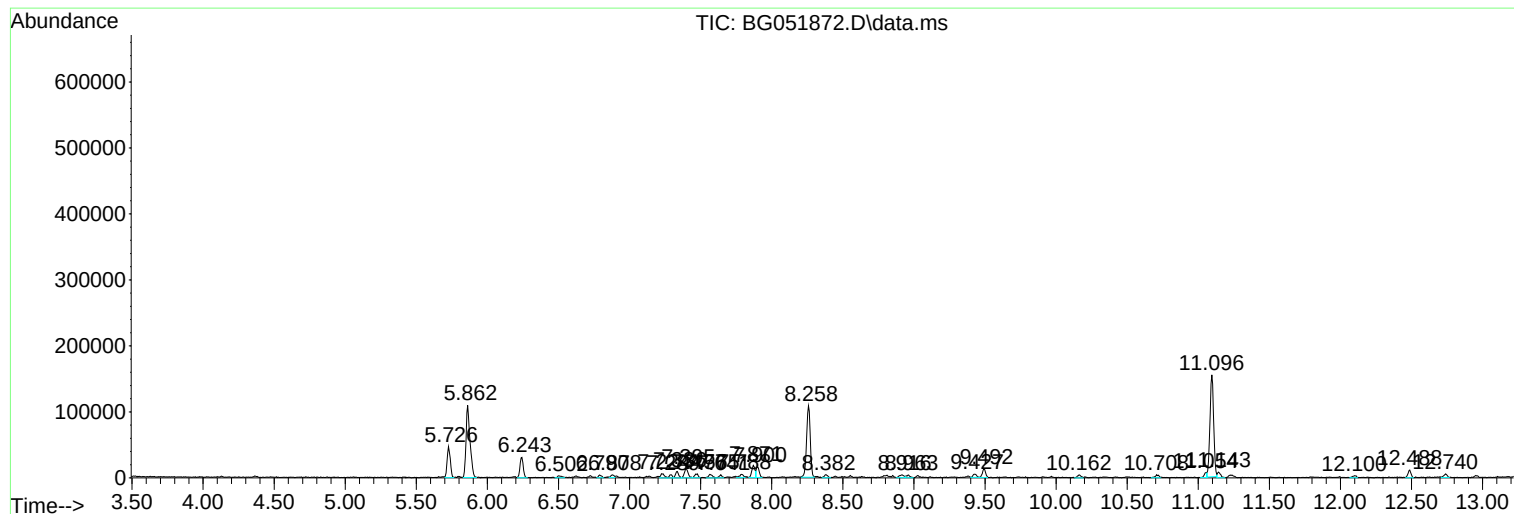
Sum of corrected areas: 6134967

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD3MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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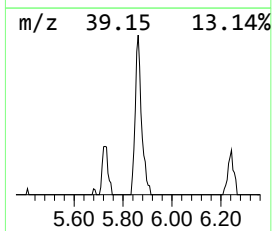
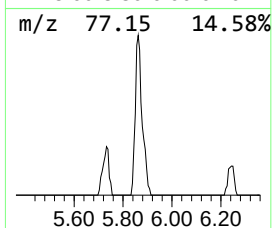
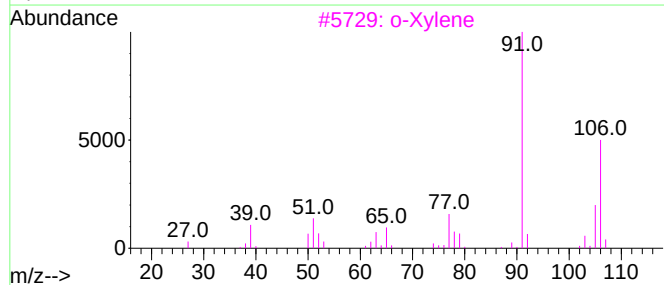
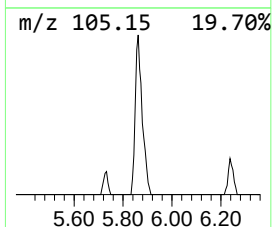
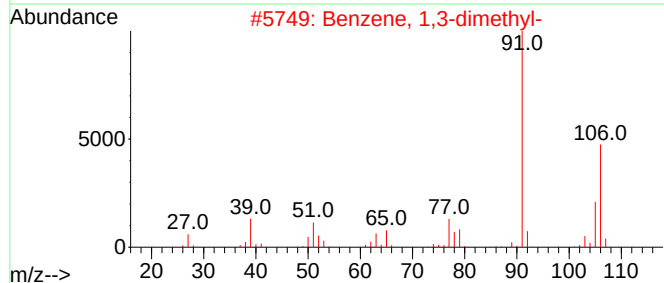
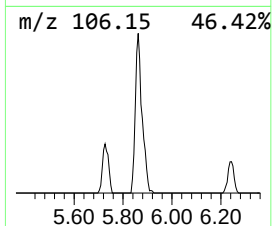
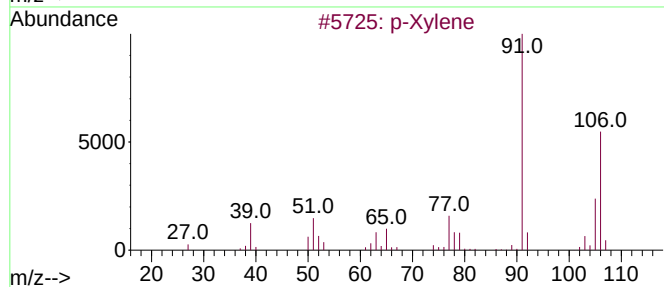
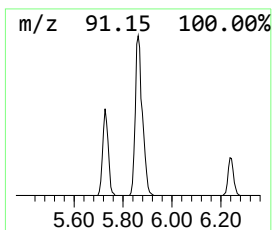
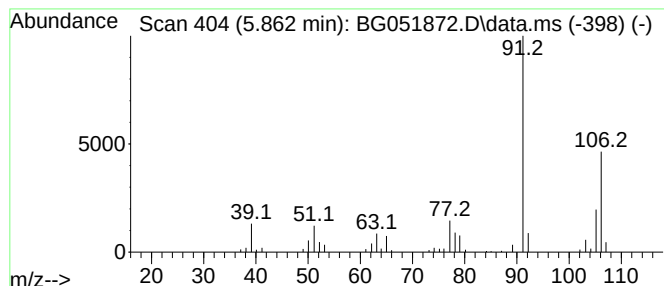
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 p-Xylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.862	21.73 ng/ul	212924	1,4-Dichlorobenzene-d4	8.258

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



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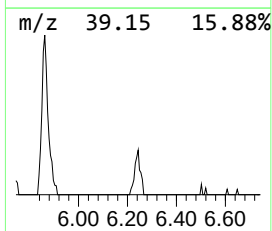
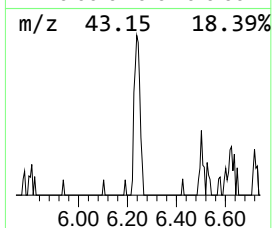
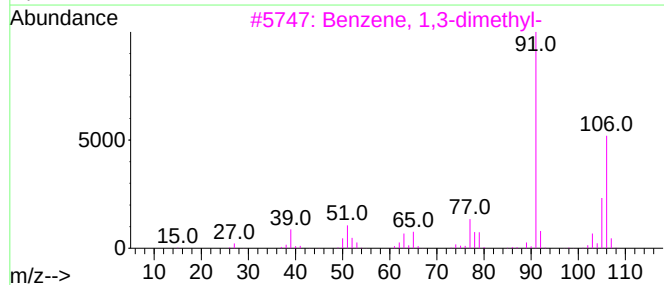
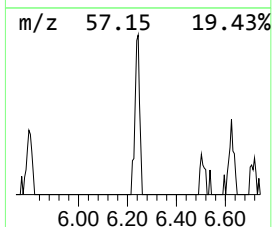
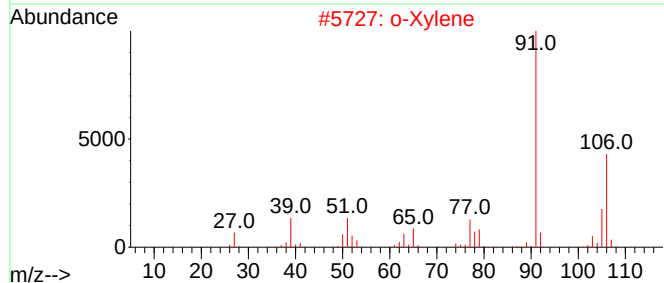
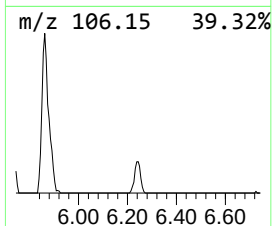
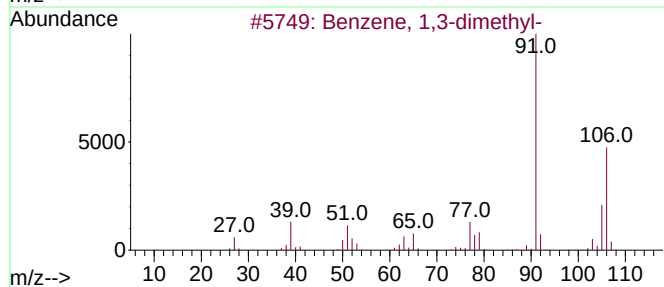
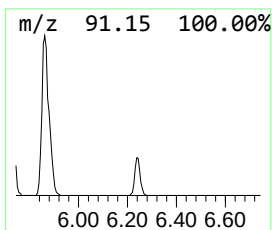
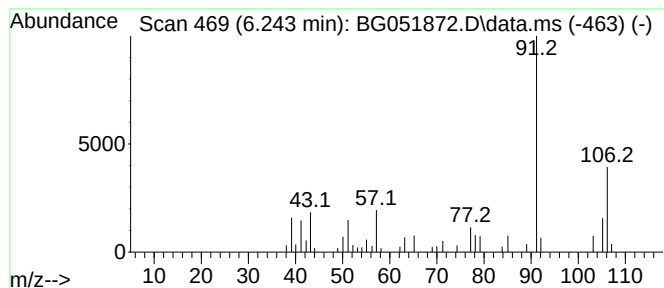
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.243	5.09 ng/ul	49898	1,4-Dichlorobenzene-d4	8.258

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
2			o-Xylene	106	C8H10	000095-47-6	91
3			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	91
4			o-Xylene	106	C8H10	000095-47-6	91
5			o-Xylene	106	C8H10	000095-47-6	91



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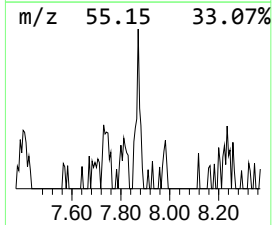
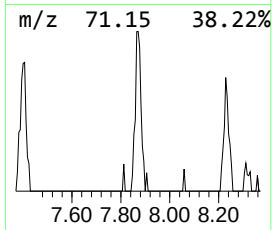
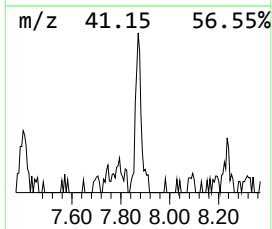
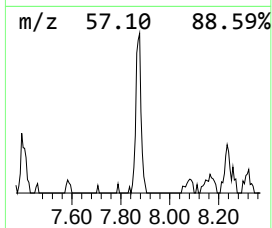
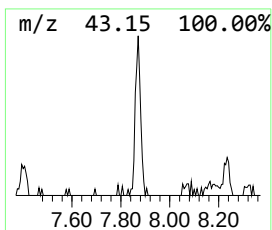
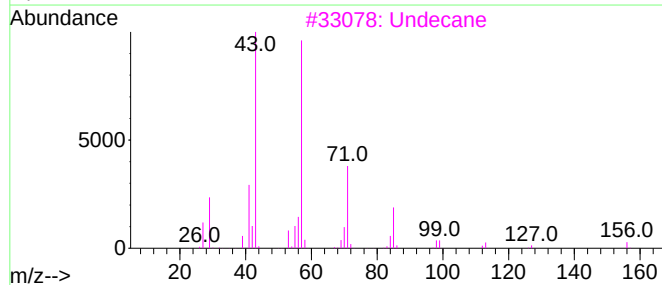
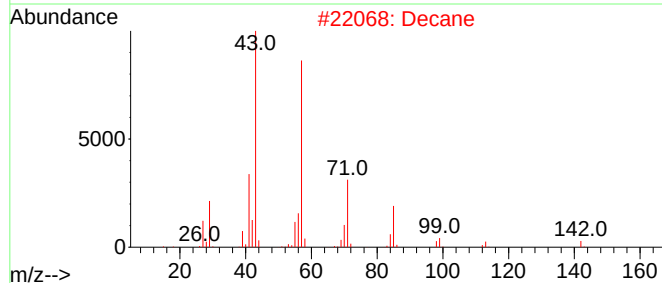
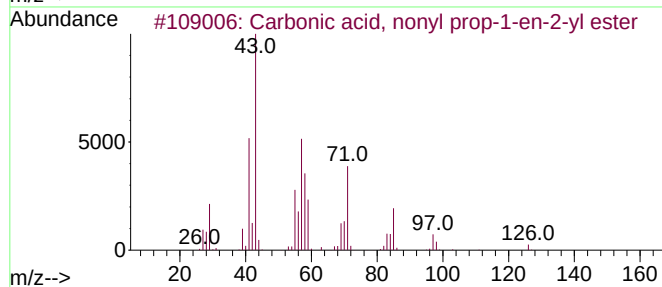
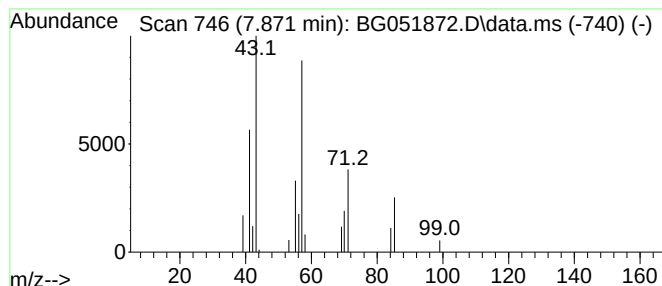
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Carbonic acid, nonyl prop-1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.871	3.09 ng/ul	30259	1,4-Dichlorobenzene-d4	8.258

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	78
2			Decane	142	C10H22	000124-18-5	72
3			Undecane	156	C11H24	001120-21-4	72
4			Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	64
5			Tridecane	184	C13H28	000629-50-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGLD3MEDL

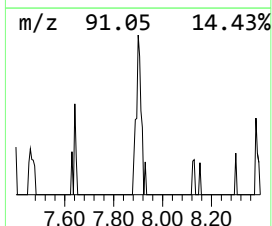
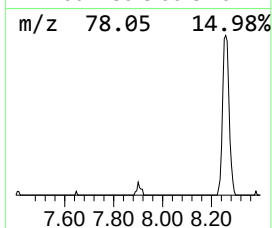
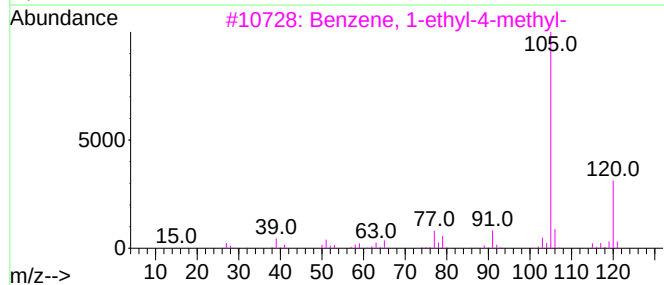
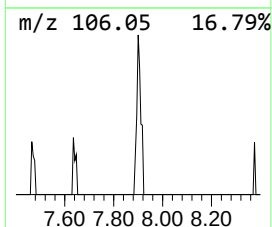
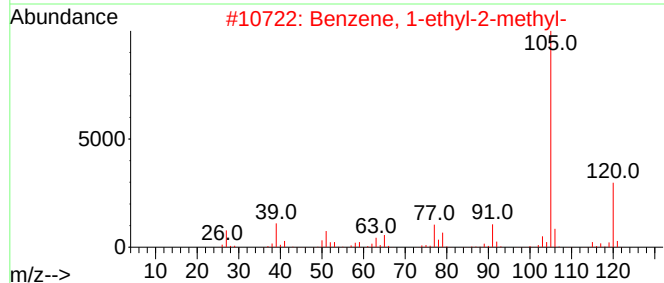
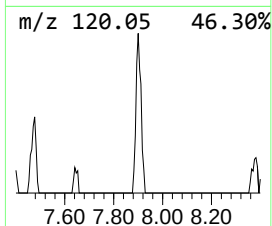
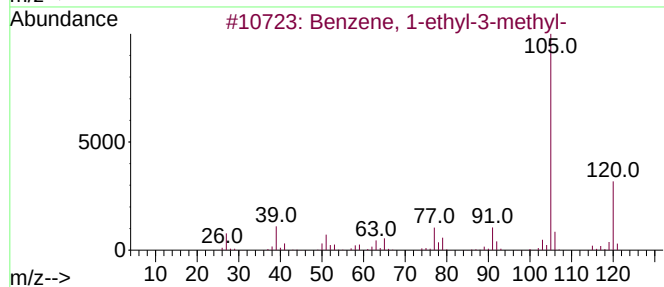
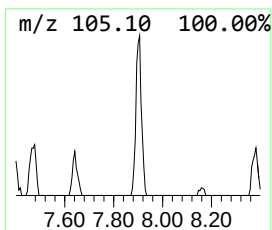
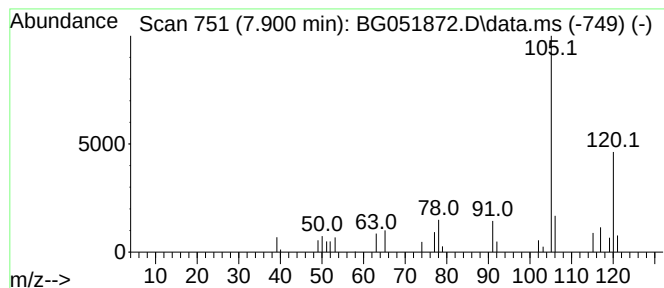
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.900	2.22 ng/ul	21723	1,4-Dichlorobenzene-d4	8.258

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	80
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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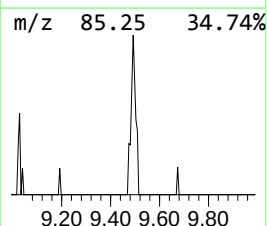
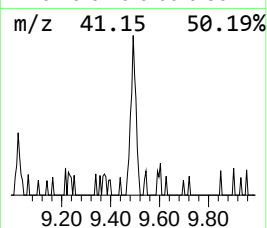
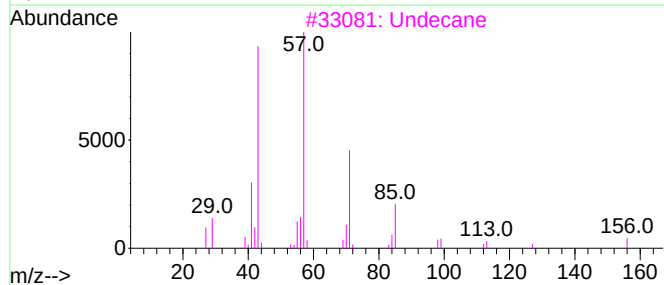
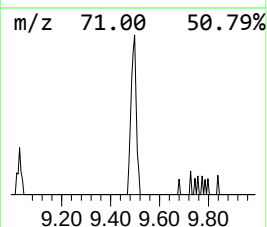
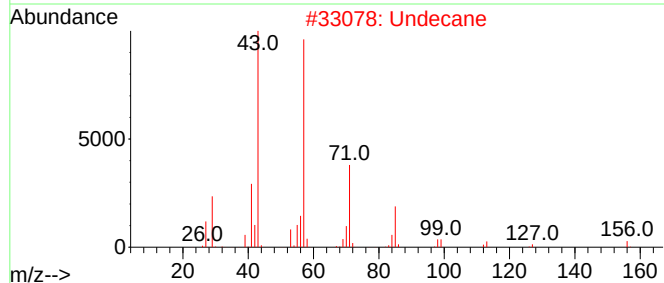
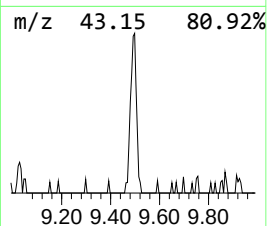
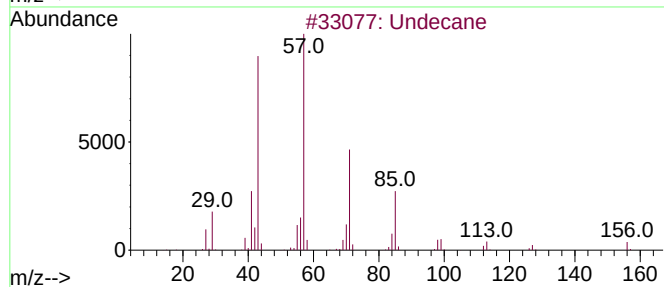
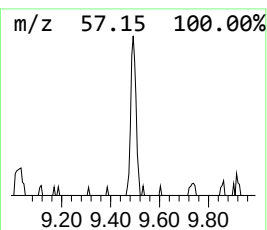
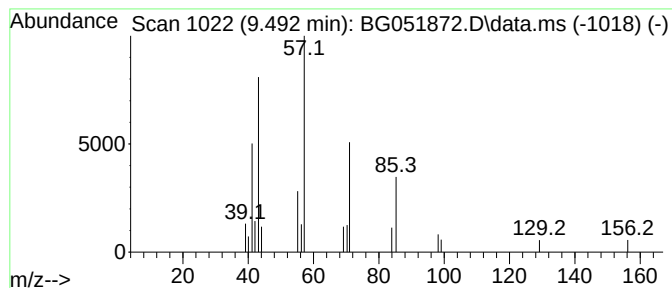
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.492	2.11 ng/ul	20635	1,4-Dichlorobenzene-d4	8.258

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	87
2	Undecane			156	C11H24	001120-21-4	83
3	Undecane			156	C11H24	001120-21-4	74
4	Undecane			156	C11H24	001120-21-4	72
5	Octacosane			394	C28H58	000630-02-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

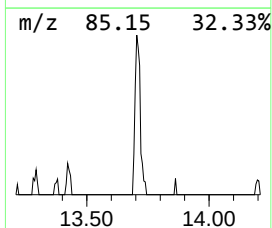
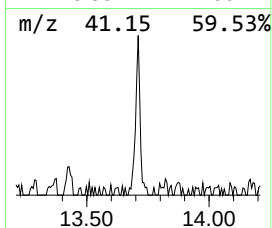
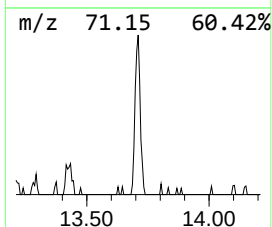
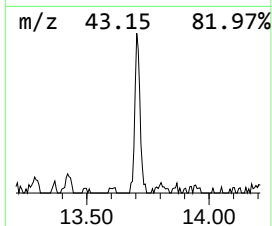
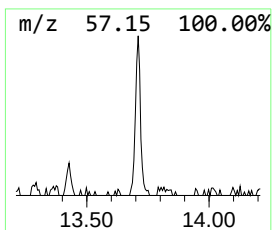
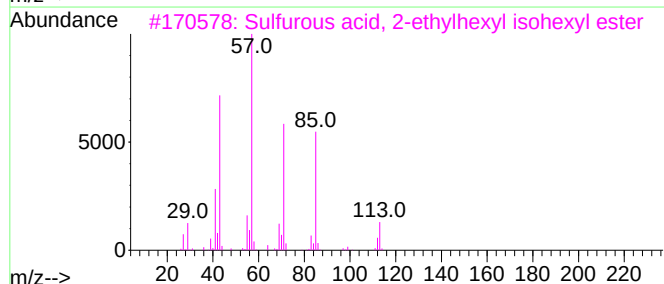
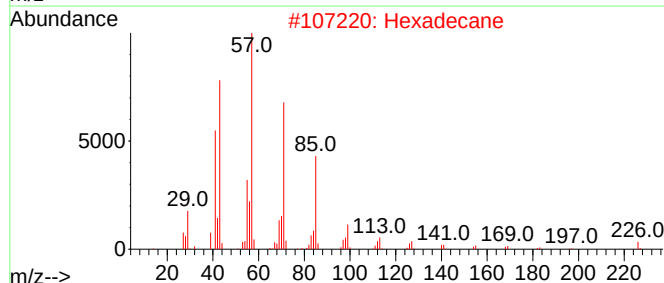
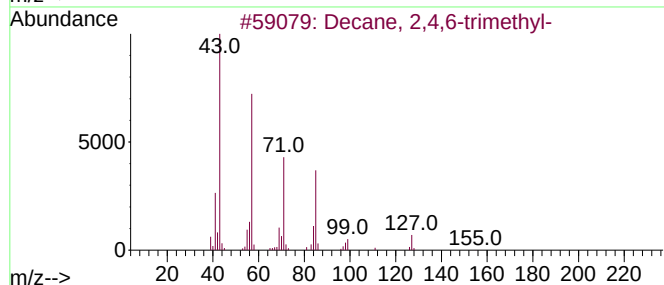
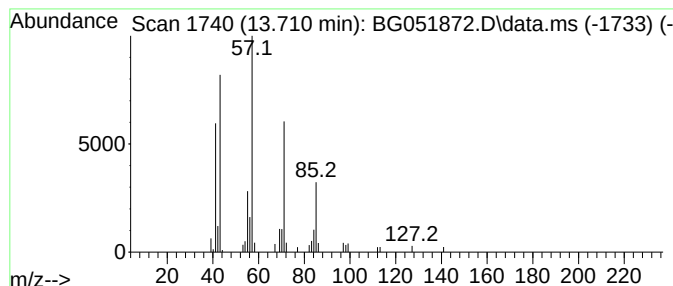
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.710	2.30 ng/ul	50714	Acenaphthene-d10	14.896	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	83
2	Hexadecane	226	C16H34	000544-76-3	72
3	Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	72
4	Hexadecane	226	C16H34	000544-76-3	72
5	Nonadecane	268	C19H40	000629-92-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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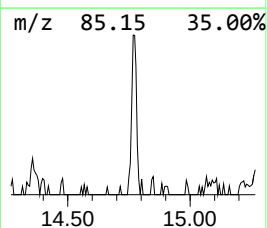
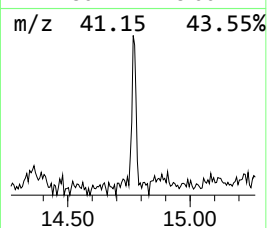
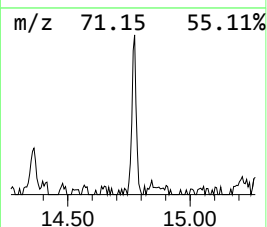
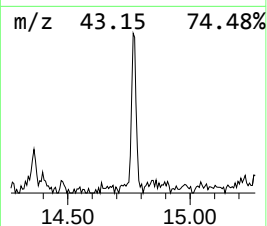
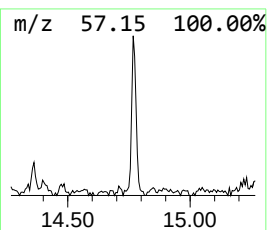
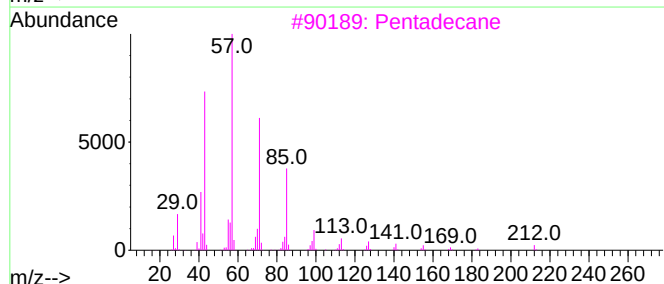
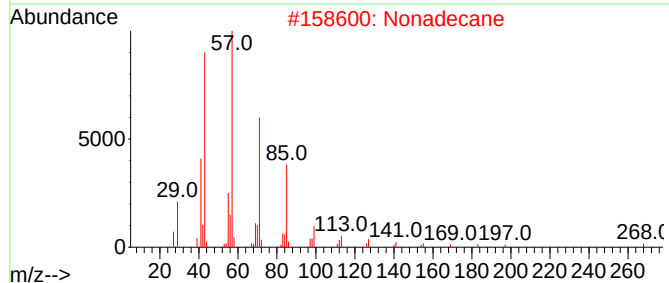
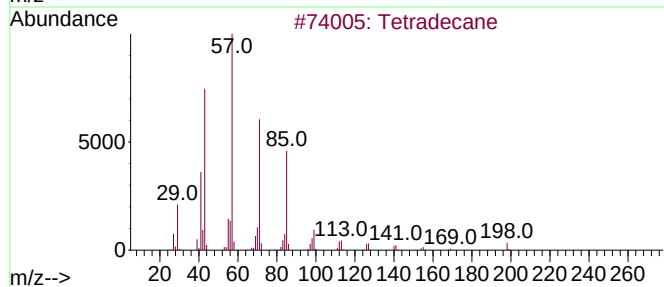
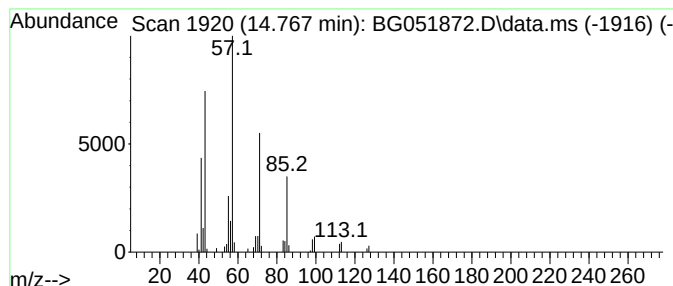
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.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.
14.767	2.46 ng/ul	54280	Acenaphthene-d10	14.896

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	90
2		Nonadecane	268	C19H40	000629-92-5	86
3		Pentadecane	212	C15H32	000629-62-9	86
4		Heptadecane	240	C17H36	000629-78-7	86
5		Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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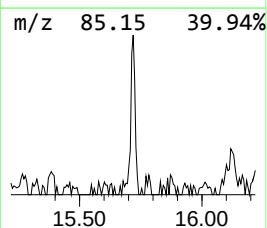
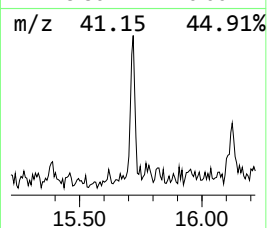
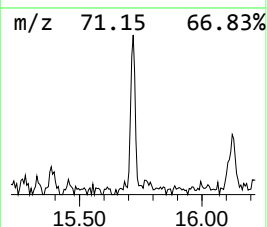
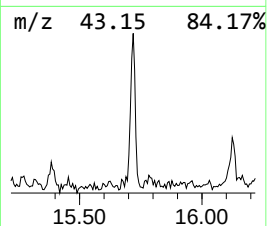
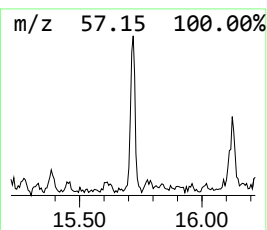
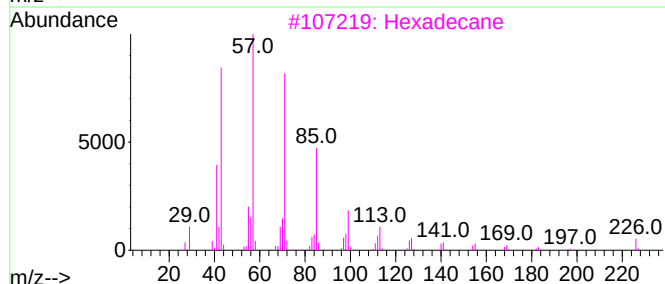
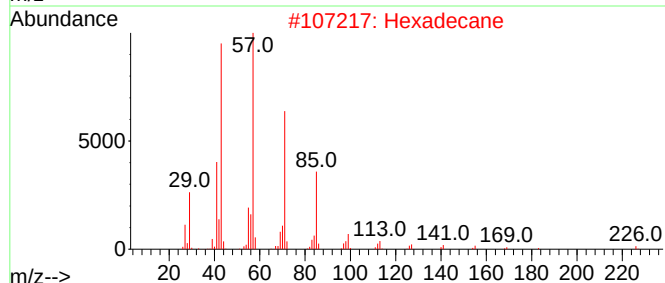
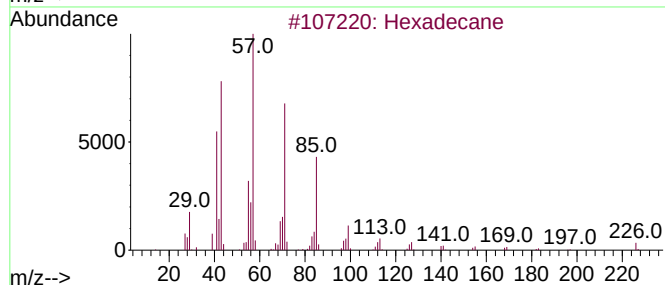
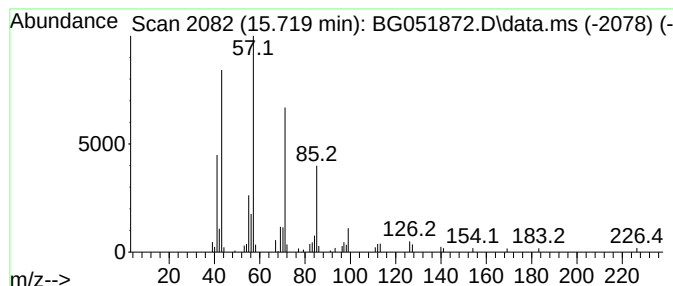
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.719	2.74 ng/ul	60252	Acenaphthene-d10	14.896

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	96
2			Hexadecane	226	C16H34	000544-76-3	96
3			Hexadecane	226	C16H34	000544-76-3	93
4			3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	90
5			Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
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Misc :
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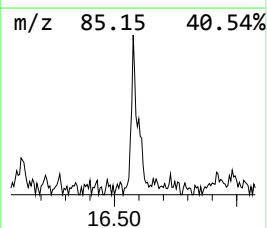
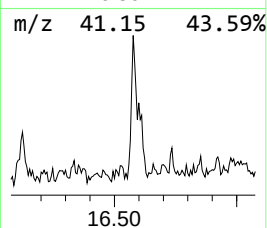
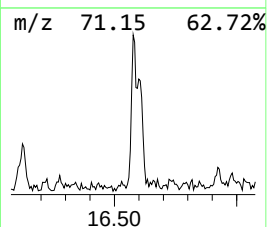
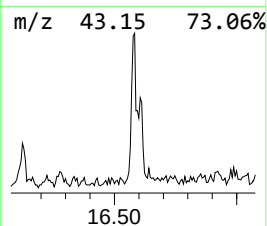
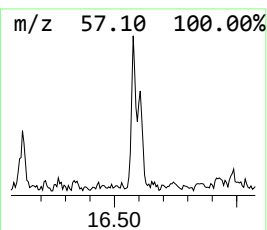
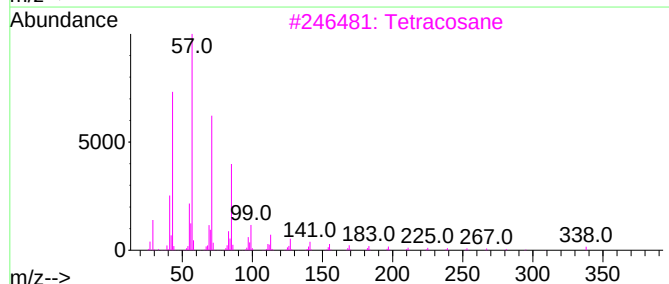
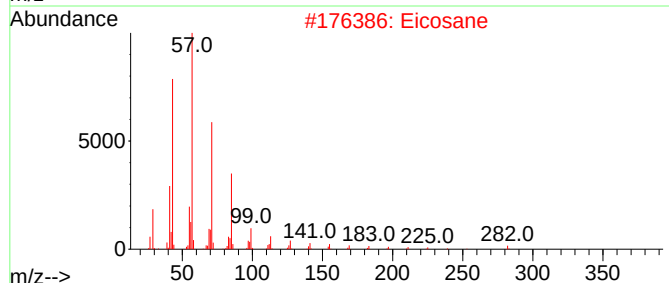
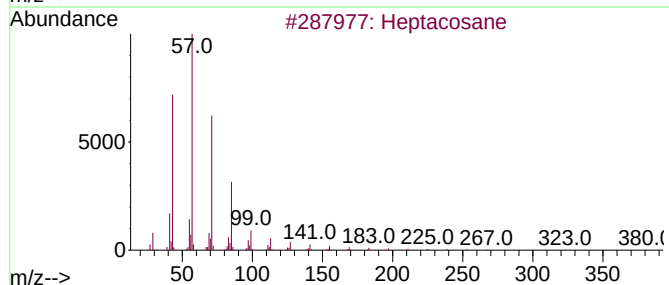
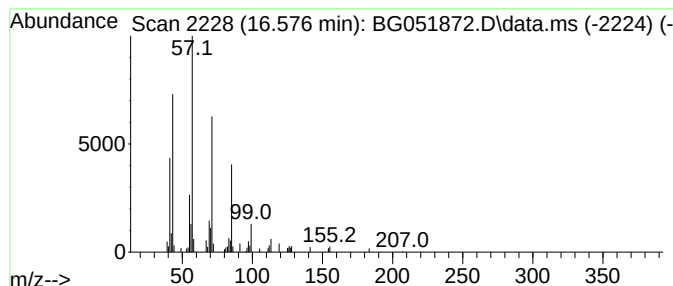
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.576	2.62 ng/ul	74763	Phenanthrene-d10	17.640

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	91
2		Eicosane	282	C20H42	000112-95-8	87
3		Tetracosane	338	C24H50	000646-31-1	86
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010422\
Data File : BG051872.D
Acq On : 4 Jan 2022 16:14
Operator : CG/JU
Sample : M5215-01MEDL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1,3-di...	6.243	5.1	ng/ul	49898	1	8.258	195975	20.0
Carbonic acid, ...	7.871	3.1	ng/ul	30259	1	8.258	195975	20.0
Benzene, 1-ethy...	7.900	2.2	ng/ul	21723	1	8.258	195975	20.0
(DEL) Alkane: S...	9.492	2.1	ng/ul	20635	1	8.258	195975	20.0
(DEL) Alkane: S...	13.710	2.3	ng/ul	50714	3	14.896	440511	20.0
(DEL) Alkane: S...	14.767	2.5	ng/ul	54280	3	14.896	440511	20.0
(DEL) Alkane: S...	15.719	2.7	ng/ul	60252	3	14.896	440511	20.0
(DEL) Alkane: S...	16.576	2.6	ng/ul	74763	4	17.640	571481	20.0