

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
 Data File : BG052390.D
 Acq On : 15 Feb 2022 12:47
 Operator : CG/JU
 Sample : N1478-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GBCW7

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-BG013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052390.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.552	7	11	16	rVB2	7457	11136	0.91%	0.042%
2	3.611	16	21	22	rBV2	2274	2178	0.18%	0.008%
3	3.910	69	72	80	rVB8	2399	3698	0.30%	0.014%
4	3.981	80	84	96	rBV	86937	145954	11.98%	0.556%
5	4.339	141	145	151	rVB4	3221	4284	0.35%	0.016%
6	4.598	184	189	193	rVB3	1544	2060	0.17%	0.008%
7	4.662	196	200	204	rBV3	3366	5352	0.44%	0.020%
8	5.279	297	305	314	rVB	28190	44162	3.62%	0.168%
9	6.384	488	493	499	rBV2	6100	8441	0.69%	0.032%
10	6.953	585	590	594	rVB2	1466	2519	0.21%	0.010%
11	7.077	605	611	616	rBV3	5255	11804	0.97%	0.045%
12	7.159	623	625	630	rBV4	1092	1961	0.16%	0.007%
13	7.218	630	635	639	rVB3	9650	14849	1.22%	0.057%
14	7.382	657	663	674	rBV2	130609	346143	28.40%	1.318%
15	7.541	683	690	699	rBV	88225	153326	12.58%	0.584%
16	7.635	699	706	717	rBV	167598	282893	23.21%	1.077%
17	7.764	722	728	736	rVB	113634	197496	16.20%	0.752%
18	8.040	771	775	778	rBV3	1105	1797	0.15%	0.007%
19	8.075	778	781	785	rBV3	1130	1560	0.13%	0.006%
20	8.234	802	808	819	rVB	82310	145715	11.96%	0.555%
21	8.469	843	848	854	rBV4	3849	6767	0.56%	0.026%
22	8.680	875	884	894	rBV	278107	480337	39.41%	1.829%
23	8.786	899	902	905	rVV3	2015	3218	0.26%	0.012%
24	8.821	905	908	913	rVB2	1141	1451	0.12%	0.006%
25	8.945	919	929	934	rBV2	151171	259378	21.28%	0.988%
26	9.004	934	939	943	rVV	287692	505339	41.46%	1.925%
27	9.045	943	946	953	rVB	212581	337023	27.65%	1.284%
28	9.121	958	959	962	rBV2	1700	1672	0.14%	0.006%
29	9.338	989	996	1002	rBV2	126142	227703	18.68%	0.867%
30	9.409	1002	1008	1014	rVB	113998	198709	16.30%	0.757%
31	9.979	1097	1105	1113	rBV	207501	356521	29.25%	1.358%
32	10.143	1125	1133	1139	rBV	101089	175617	14.41%	0.669%
33	10.225	1139	1147	1154	rBV	228703	384031	31.51%	1.463%
34	10.460	1179	1187	1193	rBV	174609	288706	23.69%	1.100%
35	10.695	1218	1227	1241	rVB2	152889	399656	32.79%	1.522%
36	11.077	1283	1292	1296	rBV	121410	222411	18.25%	0.847%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
 Title : SVOA CALIBRATION

37	11.124	1296	1300	1307	rVV	146455	259233	21.27%	0.987%
38	11.201	1307	1313	1324	rVB	83785	155261	12.74%	0.591%
39	11.342	1331	1337	1342	rVB2	7017	10916	0.90%	0.042%
40	11.412	1342	1349	1356	rBV	320572	532848	43.72%	2.029%
41	11.770	1405	1410	1416	rVB3	9049	14785	1.21%	0.056%
42	11.835	1416	1421	1425	rBV2	19174	28955	2.38%	0.110%
43	11.964	1434	1443	1459	rBV2	91179	241010	19.77%	0.918%
44	12.569	1539	1546	1555	rBV	134083	211280	17.33%	0.805%
45	12.657	1555	1561	1564	rVV3	1996	3259	0.27%	0.012%
46	12.710	1564	1570	1578	rVB	7775	10766	0.88%	0.041%
47	12.845	1587	1593	1598	rBV4	5338	12304	1.01%	0.047%
48	12.939	1600	1609	1617	rVB	371490	641578	52.64%	2.443%
49	13.080	1624	1633	1641	rVB2	304751	642226	52.69%	2.446%
50	13.398	1680	1687	1700	rBV	218393	367660	30.17%	1.400%
51	13.715	1734	1741	1742	rBV2	1123	1650	0.14%	0.006%
52	13.756	1742	1748	1757	rVB	161547	267730	21.97%	1.020%
53	14.267	1829	1835	1839	rBV	300690	443429	36.38%	1.689%
54	14.314	1839	1843	1850	rVB	178893	250587	20.56%	0.954%
55	14.573	1880	1887	1890	rBV	338309	575373	47.21%	2.191%
56	14.772	1913	1921	1931	rBV	214177	340435	27.93%	1.297%
57	14.878	1932	1939	1944	rVV	218711	326837	26.82%	1.245%
58	14.943	1944	1950	1954	rVV	382888	601934	49.39%	2.292%
59	14.978	1954	1956	1965	rVV	126188	185357	15.21%	0.706%
60	15.066	1965	1971	1979	rVV	133539	231628	19.00%	0.882%
61	15.154	1981	1986	1997	rVV	198648	284505	23.34%	1.084%
62	15.272	2000	2006	2014	rVB	204075	304186	24.96%	1.159%
63	15.724	2077	2083	2090	rBV	8452	13000	1.07%	0.050%
64	15.865	2100	2107	2111	rBV	427249	682149	55.97%	2.598%
65	15.906	2111	2114	2116	rBV	250452	349071	28.64%	1.329%
66	15.994	2123	2129	2138	rVB2	416451	758154	62.20%	2.887%
67	17.275	2341	2347	2366	rBV	463987	739790	60.70%	2.818%
68	17.621	2400	2406	2413	rBV	290449	449104	36.85%	1.710%
69	17.727	2416	2424	2426	rVV	480753	766572	62.90%	2.920%
70	17.756	2426	2429	2438	rVB	459343	673568	55.26%	2.565%
71	18.021	2467	2474	2484	rBV	516370	801793	65.78%	3.054%
72	18.162	2494	2498	2503	rBV5	2877	3229	0.26%	0.012%
73	18.232	2506	2510	2515	rVB2	13362	16972	1.39%	0.065%
74	18.561	2561	2566	2572	rVB2	24204	31434	2.58%	0.120%
75	18.673	2581	2585	2590	rVB2	9399	12037	0.99%	0.046%
76	18.779	2598	2603	2608	rBV	33949	49648	4.07%	0.189%

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77	18.826	2608	2611	2617	rVB2	19647	23366	1.92%	0.089%
78	18.984	2635	2638	2641	rBV3	1452	1782	0.15%	0.007%
79	19.031	2641	2646	2651	rVV2	45237	60866	4.99%	0.232%
80	19.119	2656	2661	2664	rBV4	2735	3863	0.32%	0.015%
81	19.260	2680	2685	2689	rVB3	2966	4017	0.33%	0.015%
82	19.566	2732	2737	2740	rBV4	7605	10068	0.83%	0.038%
83	19.601	2740	2743	2746	rVV3	7296	8657	0.71%	0.033%
84	19.666	2746	2754	2762	rVB	391212	564686	46.33%	2.151%
85	19.736	2762	2766	2769	rBV4	3656	4761	0.39%	0.018%
86	19.771	2771	2772	2776	rVB4	3156	3079	0.25%	0.012%
87	20.000	2805	2811	2817	rBV	678535	966052	79.26%	3.679%
88	20.136	2830	2834	2839	rVB3	4595	6876	0.56%	0.026%
89	20.893	2958	2963	2969	rVB	373108	476242	39.07%	1.814%
90	21.792	3110	3116	3125	rVB	385577	545349	44.74%	2.077%
91	21.939	3134	3141	3145	rBV	309988	544549	44.68%	2.074%
92	21.986	3145	3149	3157	rVB	581907	989577	81.19%	3.769%
93	22.767	3278	3282	3286	rBV6	4831	9642	0.79%	0.037%
94	23.038	3321	3328	3336	rVB	86291	170880	14.02%	0.651%
95	24.371	3545	3555	3566	rBV	469225	1218811	100.00%	4.642%
96	25.170	3680	3691	3698	rBV2	306350	949451	77.90%	3.616%
97	25.252	3698	3705	3717	rVB	378904	1069101	87.72%	4.072%
98	25.411	3722	3732	3744	rVB2	180905	551523	45.25%	2.100%
99	26.674	3939	3947	3954	rVB9	14031	43416	3.56%	0.165%
100	29.411	4400	4413	4432	rVB3	99805	502148	41.20%	1.912%

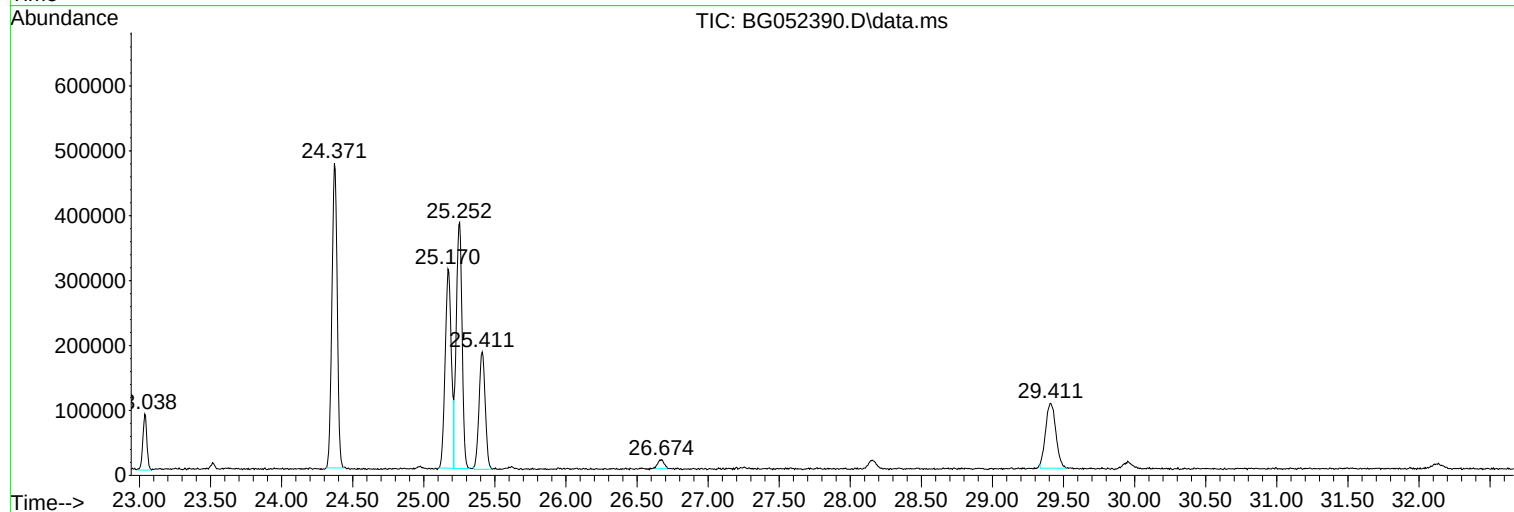
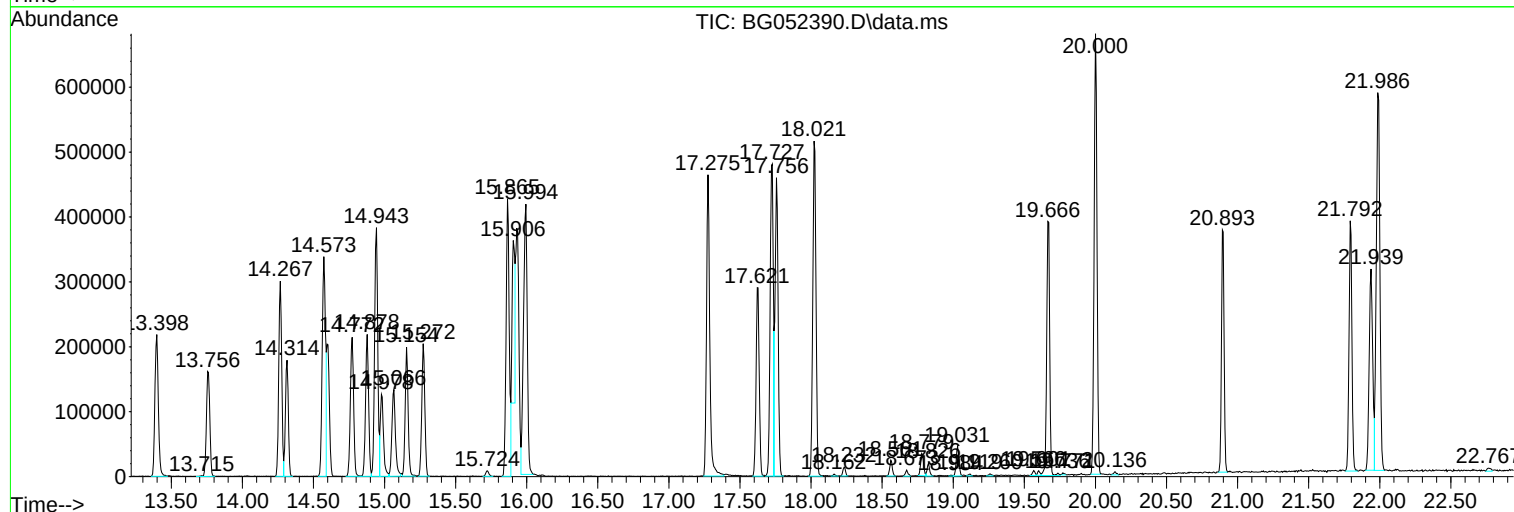
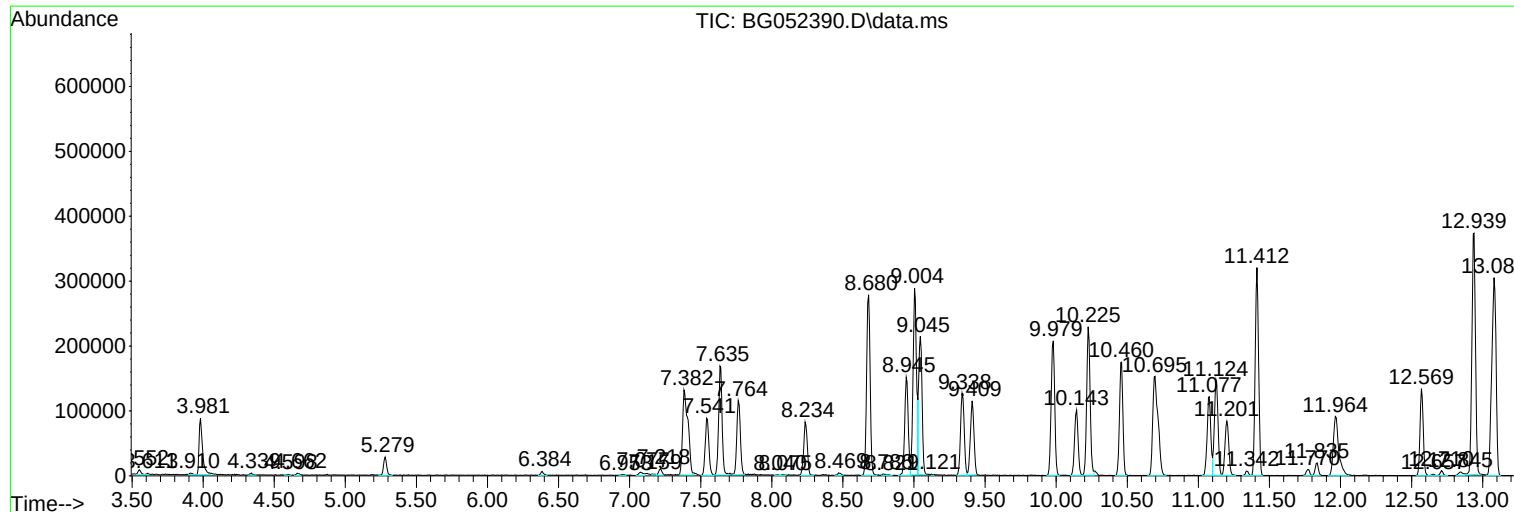
Sum of corrected areas: 26256882

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ALS Vial : 4 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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



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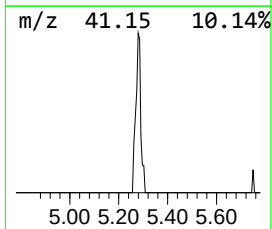
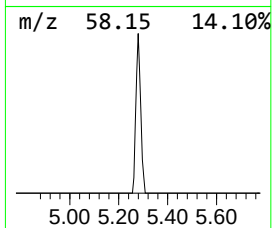
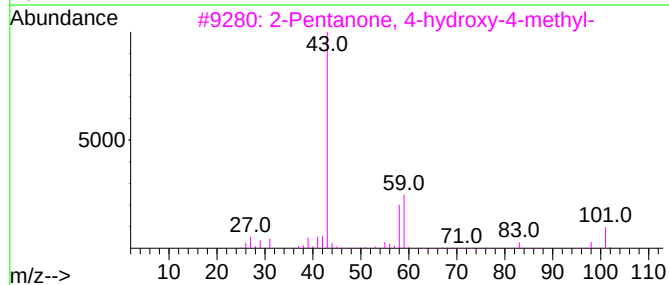
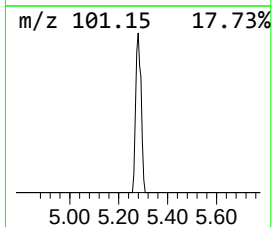
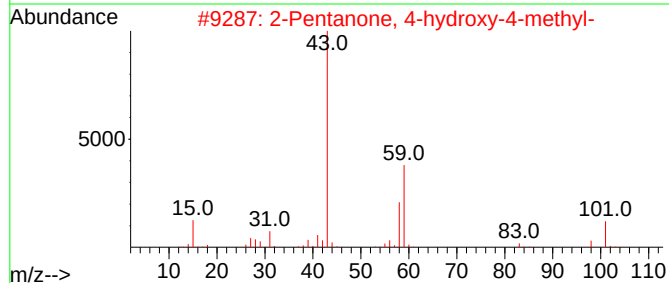
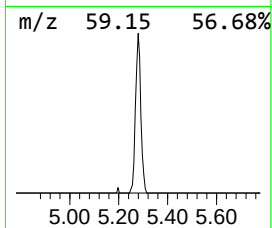
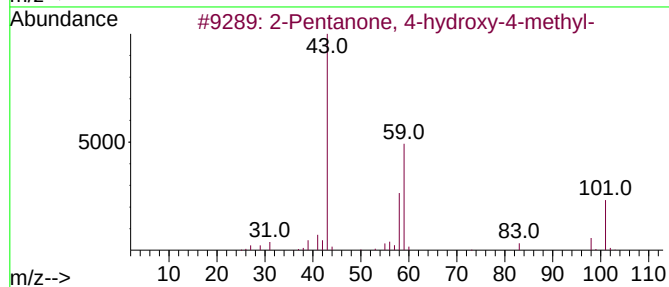
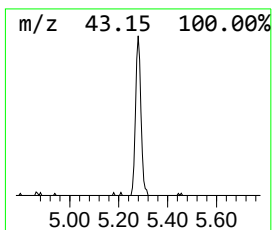
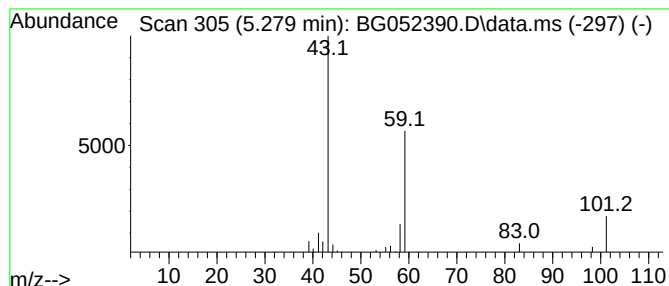
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.279	6.06 ng/ul	44162	1,4-Dichlorobenzene-d4	8.240

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39		
5	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	36		



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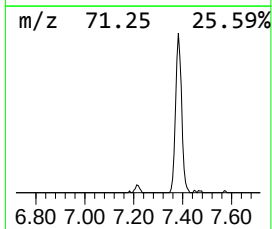
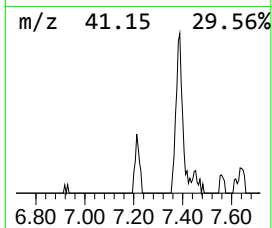
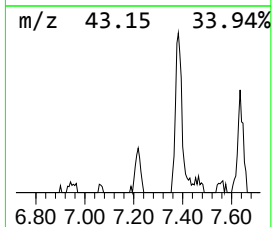
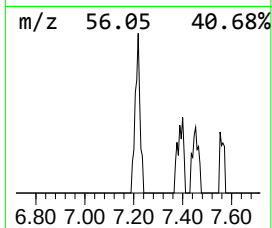
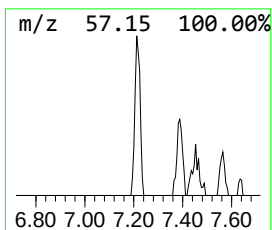
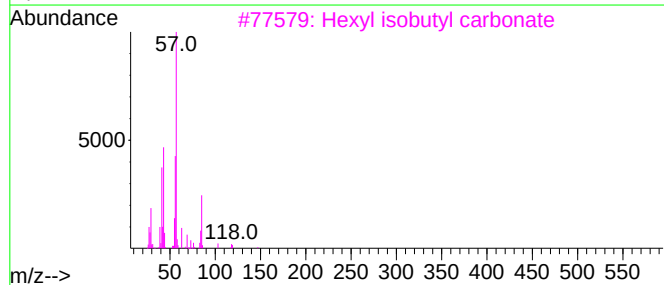
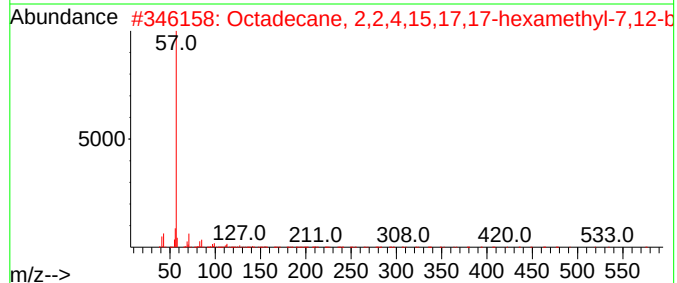
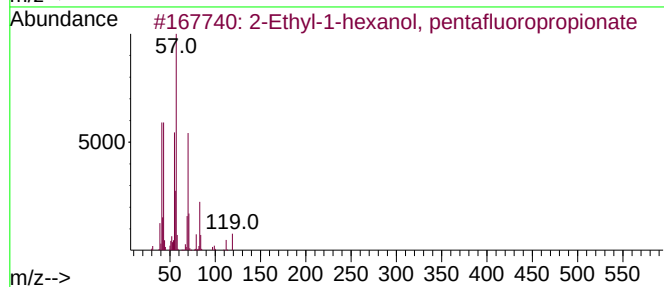
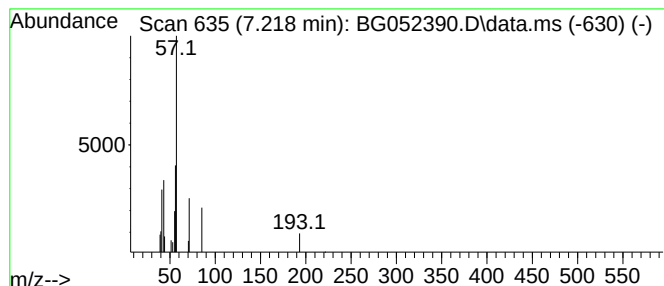
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.218	2.04 ng/ul	14849	1,4-Dichlorobenzene-d4	8.240

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethyl-1-hexanol, pentafluoropr...	276	C11H17F5O2	959012-82-9	36
2			Octadecane, 2,2,4,15,17,17-hexam...	591	C42H86	055470-97-8	33
3			Hexyl isobutyl carbonate	202	C11H22O3	959067-98-2	28
4			Oxalic acid, 6-ethyloct-3-yl eth...	258	C14H26O4	1000309-33-9	12
5			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	9



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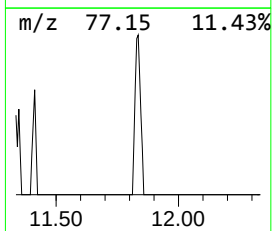
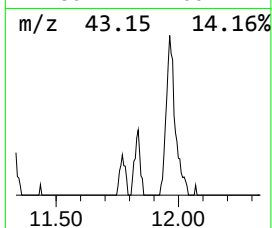
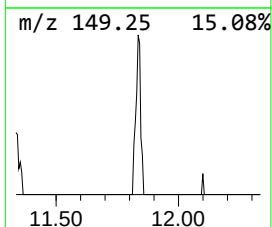
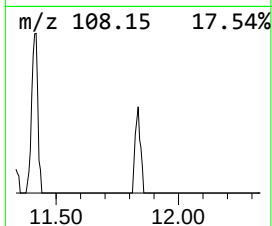
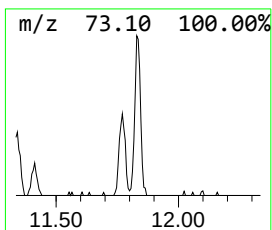
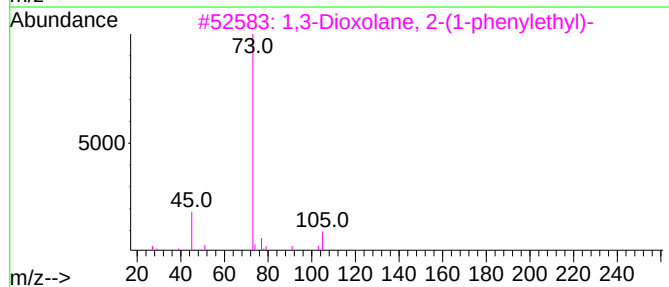
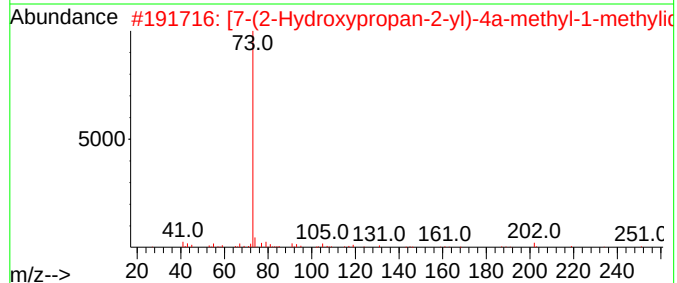
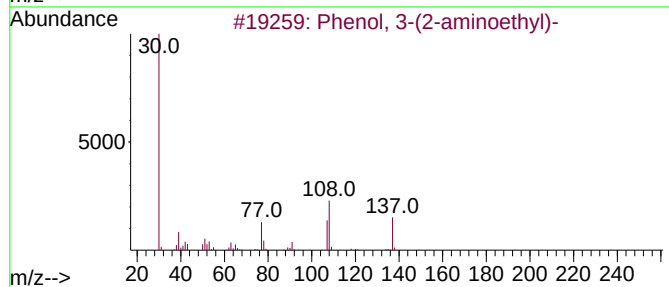
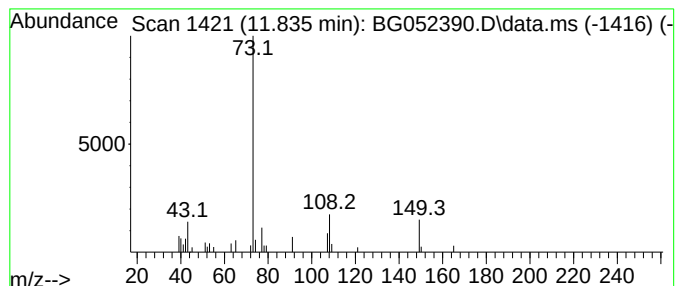
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.835	2.60 ng/ul	28955	Naphthalene-d8	11.077

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3-(2-aminoethyl)-	137	C8H11NO	000588-05-6	16
2			[7-(2-Hydroxypropan-2-yl)-4a-met...	294	C18H30O3	1010502-54-4	9
3			1,3-Dioxolane, 2-(1-phenylethyl)-	178	C11H14O2	004362-22-5	9
4			Phenol, 2-(2-aminoethyl)-	137	C8H11NO	002039-66-9	9
5			Phenol, 4-(2-aminoethyl)-	137	C8H11NO	000051-67-2	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
Operator : CG/JU
Sample : N1478-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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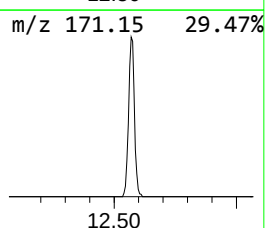
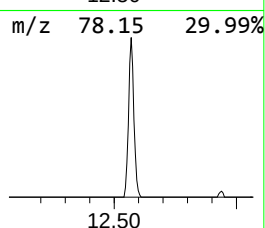
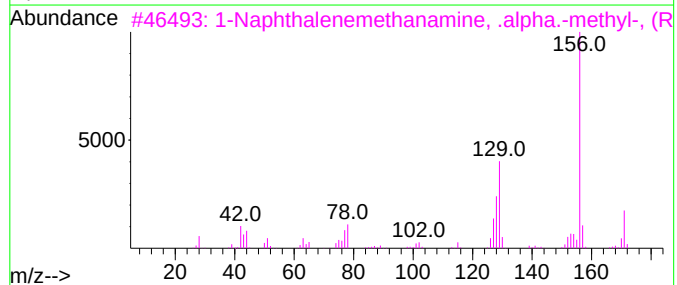
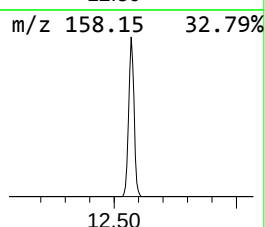
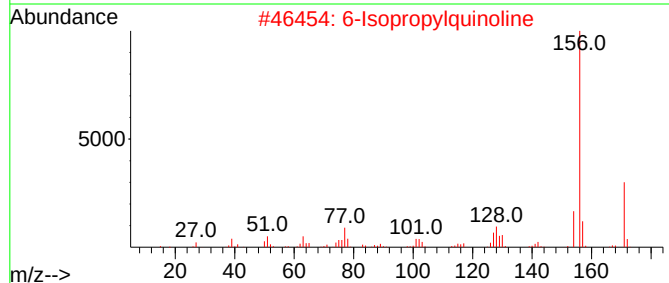
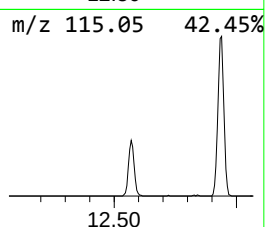
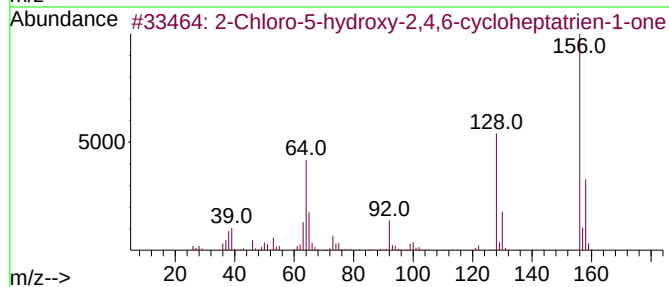
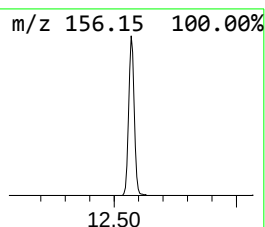
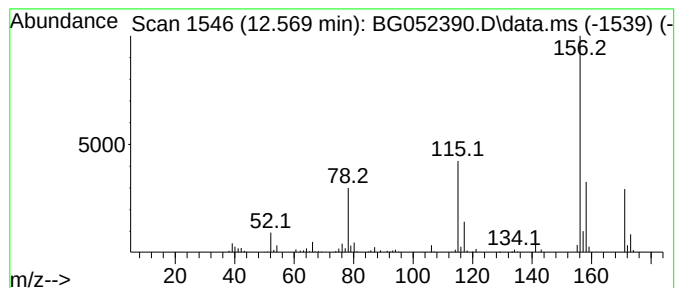
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.569	19.00 ng/ul	211280	Naphthalene-d8	11.077

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2			6-Isopropylquinoline	171	C12H13N	000135-79-5	43
3			1-Naphthalenemethanamine, .alpha...	171	C12H13N	003886-70-2	43
4			Quinoline, 8-propyl-	171	C12H13N	007661-53-2	43
5			Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
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Sample : N1478-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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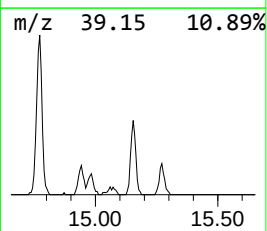
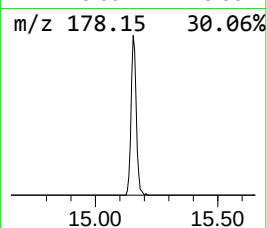
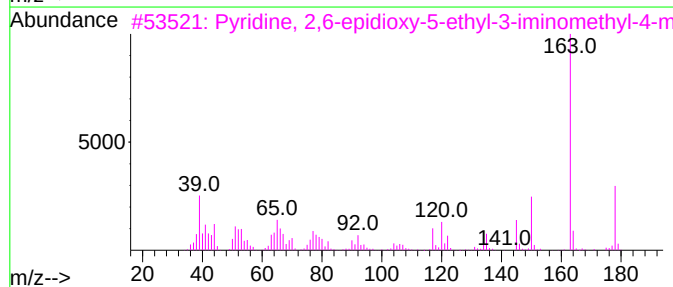
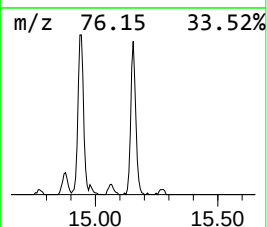
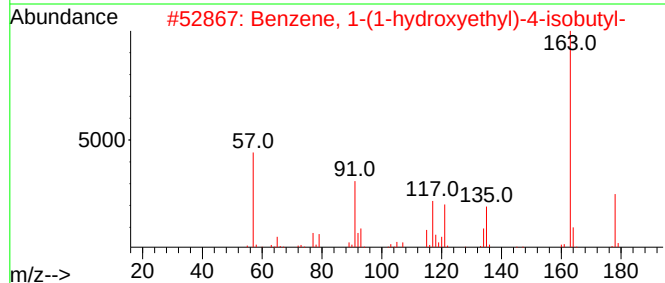
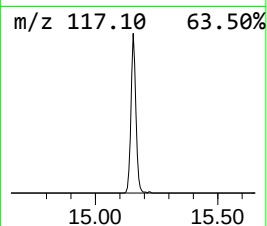
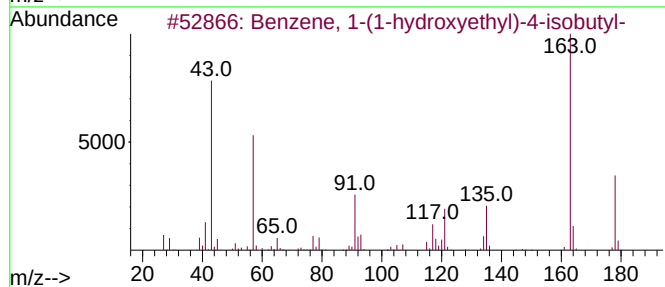
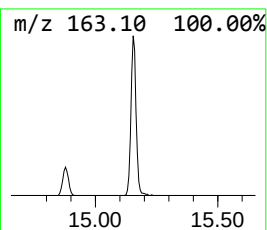
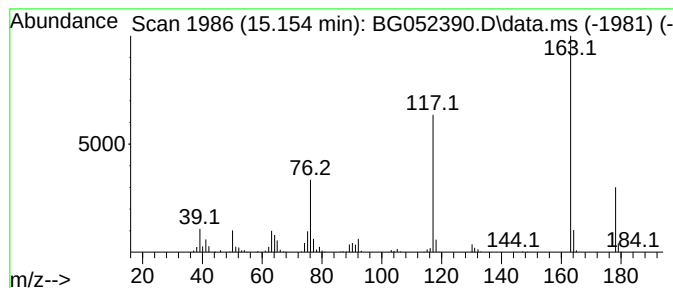
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-(1-hydroxyethyl)... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.154	17.41 ng/ul	284505	Acenaphthene-d10	14.878

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	58
2			Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	53
3			Pyridine, 2,6-epidioxo-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	50
4			Propofol	178	C12H18O	002078-54-8	50
5			2,3,4,5,6-Pentamethylphenol, met...	178	C12H18O	014804-37-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
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Sample : N1478-09
Misc :
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Instrument :
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ClientSampleId :
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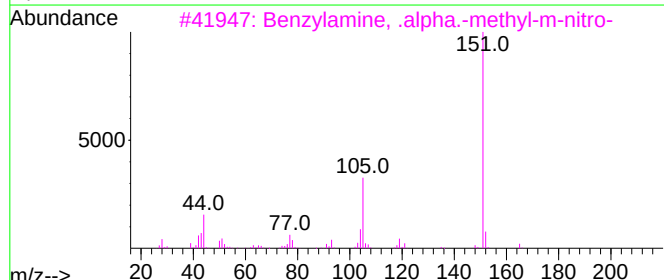
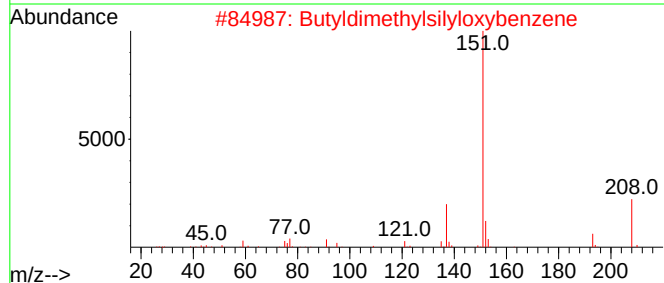
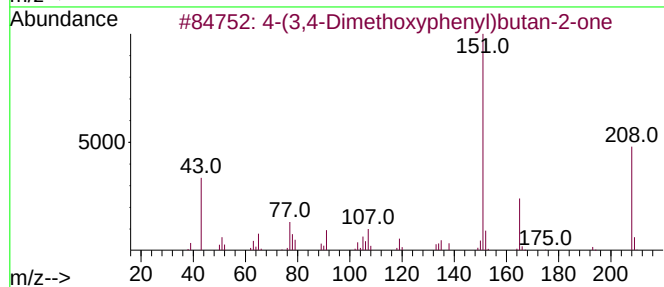
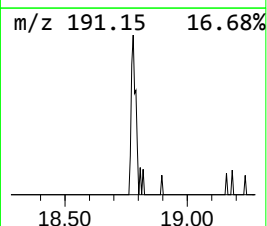
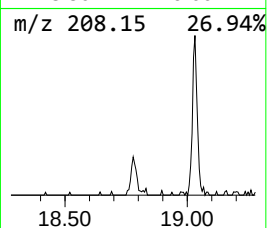
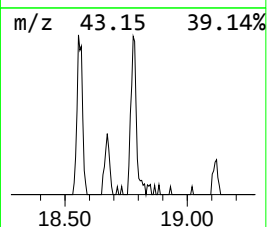
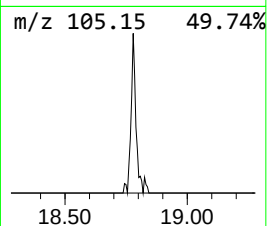
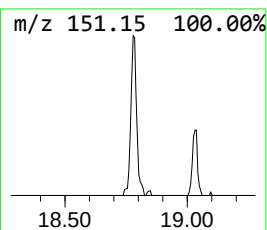
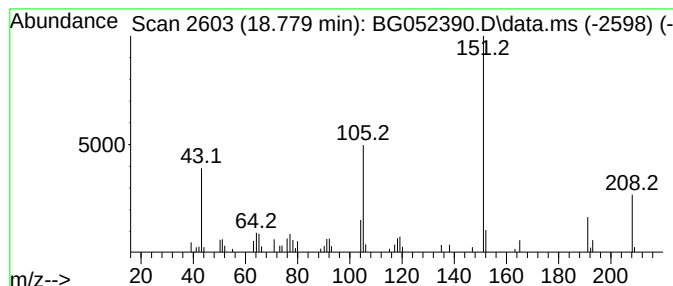
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 4-(3,4-Dimethoxyphenyl)buta... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.779	2.21 ng/ul	49648	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(3,4-Dimethoxyphenyl)butan-2-one	208	C12H16O3	006302-60-9	64
2			Butyldimethylsilyloxybenzene	208	C12H20OSi	1000299-06-1	64
3			Benzylamine, .alpha.-methyl-m-ni...	166	C8H10N2O2	090271-37-7	50
4			Ethanol, 2-[(2-nitrophenyl)amino]-	182	C8H10N2O3	004926-55-0	47
5			1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
Operator : CG/JU
Sample : N1478-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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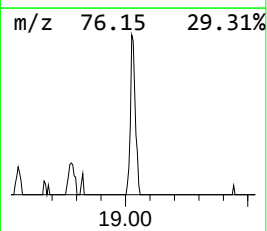
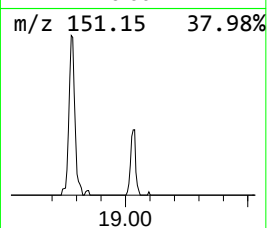
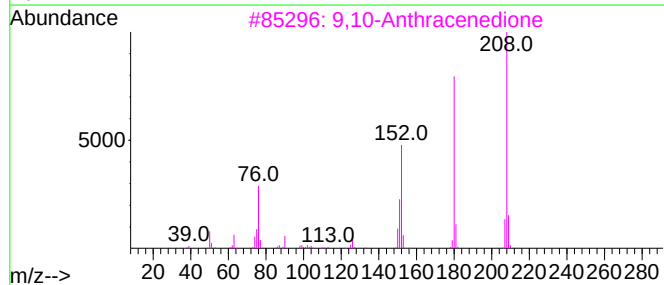
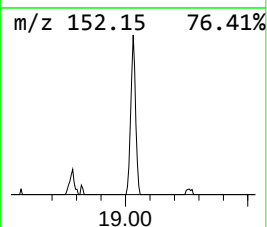
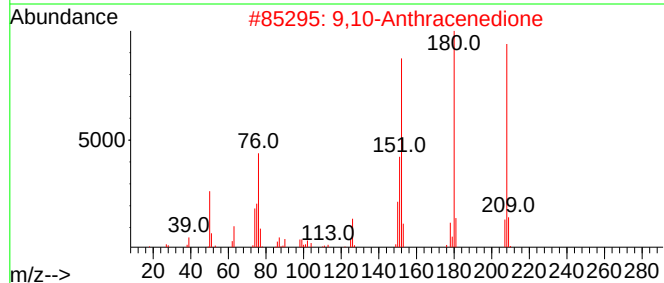
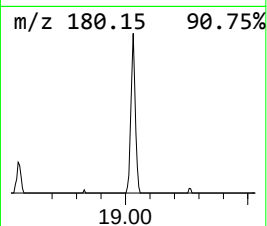
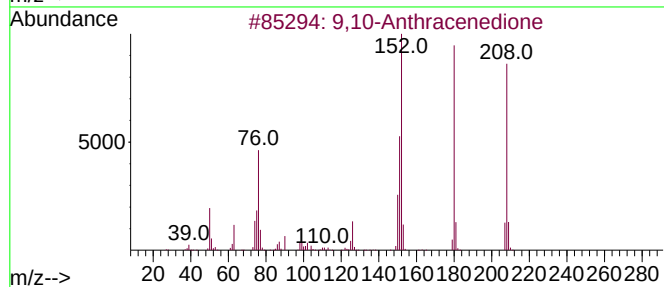
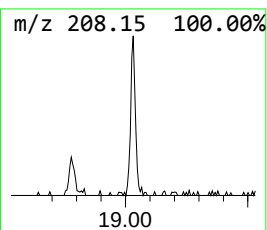
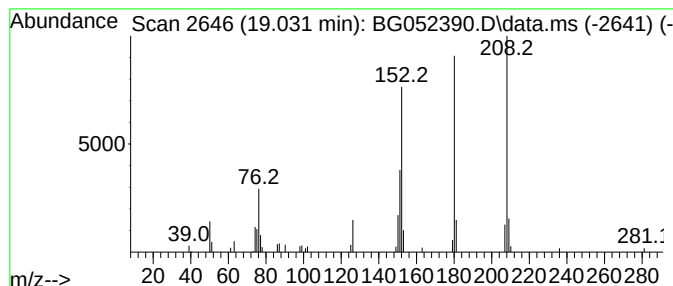
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.031	2.71 ng/ul	60866	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	97
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
4	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	91
5	1H-Phenalen-1-one			180	C13H8O	000548-39-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
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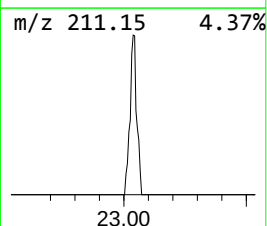
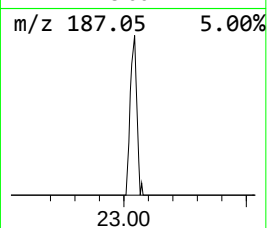
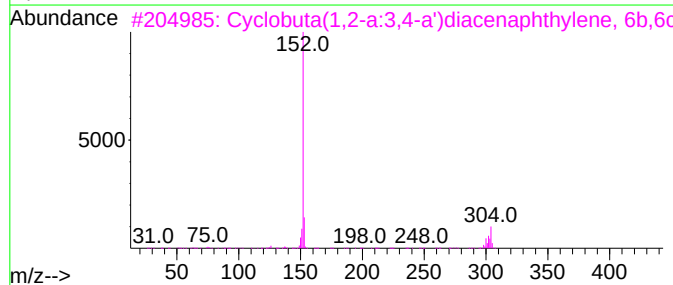
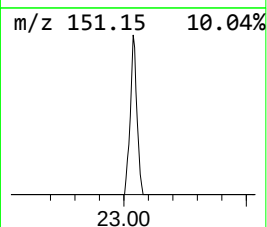
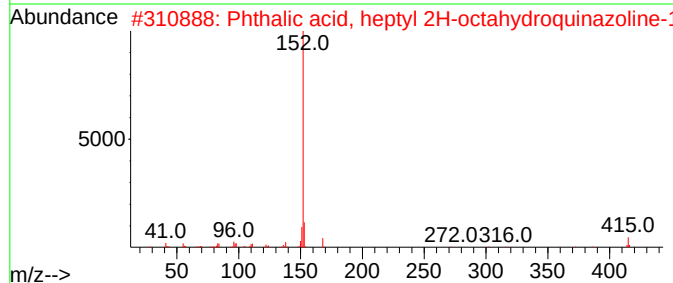
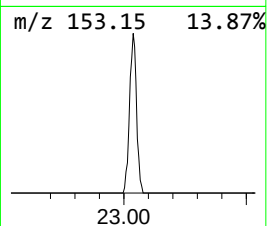
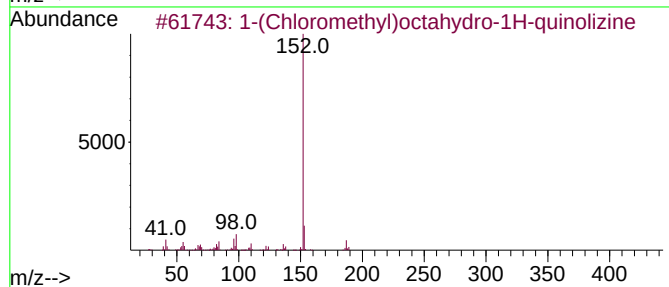
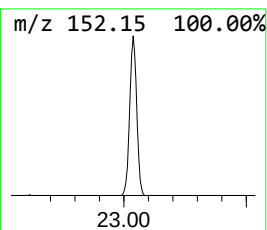
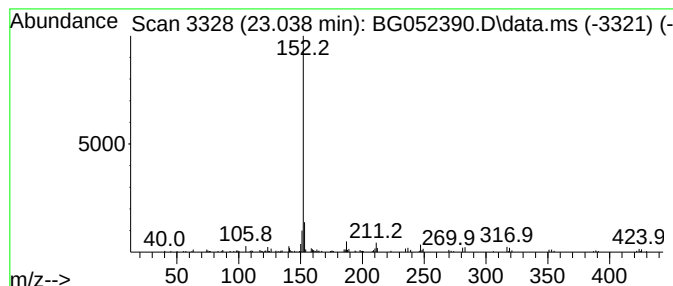
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-(Chloromethyl)octahydro-1... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.038	6.28 ng/ul	170880	Chrysene-d12	21.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(Chloromethyl)octahydro-1H-qui...	187	C10H18ClN	091049-90-0	59
2			Phthalic acid, heptyl 2H-octahyd...	415	C25H37NO4	1000322-80-8	50
3			Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	50
4			Phthalic acid, isohexyl 2H-octah...	401	C24H35NO4	1000322-80-6	50
5			2-Hydroxy-5-methoxybenzaldehyde,...	208	C12H16O3	1000395-40-1	39



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021522\
Data File : BG052390.D
Acq On : 15 Feb 2022 12:47
Operator : CG/JU
Sample : N1478-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.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
2-Pentanone, 4-...	5.279	6.1	ng/ul	44162	1	8.240	145715	20.0
unknown-01	7.218	2.0	ng/ul	14849	1	8.240	145715	20.0
unknown-02	11.835	2.6	ng/ul	28955	2	11.077	222411	20.0
unknown-03	12.569	19.0	ng/ul	211280	2	11.077	222411	20.0
Benzene, 1-(1-h...	15.154	17.4	ng/ul	284505	3	14.878	326837	20.0
4-(3,4-Dimethox...	18.779	2.2	ng/ul	49648	4	17.621	449104	20.0
9,10-Anthracene...	19.031	2.7	ng/ul	60866	4	17.621	449104	20.0
1-(Chloromethyl...	23.038	6.3	ng/ul	170880	5	21.939	544549	20.0