

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
 Data File : BG059601.D
 Acq On : 2 Nov 2023 13:33
 Operator : MA/JU
 Sample : 04996-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4A86

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BG102723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG059601.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.971	94	99	106	rBV6	30975	77549	0.11%	0.042%
2	4.083	113	118	127	rBV	224150	379923	0.55%	0.205%
3	4.441	168	179	197	rVB	35916211	68674596	100.00%	37.077%
4	4.758	226	233	244	rBV	709626	1042560	1.52%	0.563%
5	5.393	334	341	348	rVB	1503365	2256406	3.29%	1.218%
6	6.486	520	527	535	rVB	716229	1144496	1.67%	0.618%
7	7.467	685	694	695	rBV	424661	617209	0.90%	0.333%
8	7.496	695	699	707	rVB	566010	995178	1.45%	0.537%
9	7.673	721	729	733	rBV	254131	449909	0.66%	0.243%
10	7.714	733	736	740	rVV3	81241	155151	0.23%	0.084%
11	7.767	740	745	750	rVB	461204	772027	1.12%	0.417%
12	7.866	756	762	772	rVB	389445	682823	0.99%	0.369%
13	8.354	837	845	852	rBV	295002	517056	0.75%	0.279%
14	9.036	954	961	970	rBV	535408	911827	1.33%	0.492%
15	9.153	975	981	988	rBV2	243923	409916	0.60%	0.221%
16	9.547	1042	1048	1053	rBV2	338369	634064	0.92%	0.342%
17	9.594	1053	1056	1066	rVB2	130389	226556	0.33%	0.122%
18	9.817	1087	1094	1103	rVB	985672	1598928	2.33%	0.863%
19	10.269	1163	1171	1174	rBV2	311615	597144	0.87%	0.322%
20	10.305	1174	1177	1189	rVB2	251506	407768	0.59%	0.220%
21	10.798	1252	1261	1273	rBV3	756458	2292681	3.34%	1.238%
22	11.204	1322	1330	1334	rBV	575140	1077417	1.57%	0.582%
23	11.251	1334	1338	1346	rVB	434256	760701	1.11%	0.411%
24	11.345	1346	1354	1362	rBV2	349807	641429	0.93%	0.346%
25	11.480	1370	1377	1384	rBV	643756	1125972	1.64%	0.608%
26	12.678	1574	1581	1587	rBV	624890	1000321	1.46%	0.540%
27	12.831	1600	1607	1616	rBV	1103857	1773262	2.58%	0.957%
28	12.931	1619	1624	1635	rVV5	109105	252316	0.37%	0.136%
29	13.049	1637	1644	1651	rVV	1674548	2736155	3.98%	1.477%
30	13.125	1651	1657	1662	rVV2	953852	1565339	2.28%	0.845%
31	13.172	1662	1665	1671	rVB2	379660	580303	0.85%	0.313%
32	13.413	1697	1706	1712	rBV2	1237484	1964370	2.86%	1.061%
33	13.818	1768	1775	1780	rBV	2351681	3744761	5.45%	2.022%
34	13.871	1780	1784	1791	rVB	496332	809659	1.18%	0.437%
35	14.376	1863	1870	1880	rBV	1426643	2055250	2.99%	1.110%
36	14.564	1895	1902	1909	rBV	1142641	1667130	2.43%	0.900%

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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_G\Methods\SFAM-EPA-BG102723.MA.M
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37	14.688	1916	1923	1926	rBV2	1516300	2700708	3.93%	1.458%
38	14.982	1967	1973	1979	rBV	941912	1448634	2.11%	0.782%
39	15.046	1979	1984	1988	rVV	669923	1025926	1.49%	0.554%
40	15.099	1988	1993	1997	rVV	947722	1396994	2.03%	0.754%
41	15.152	1997	2002	2012	rVB2	793417	1281979	1.87%	0.692%
42	15.252	2012	2019	2029	rBV2	2271659	5535707	8.06%	2.989%
43	15.381	2035	2041	2047	rVB	685953	1021291	1.49%	0.551%
44	15.969	2135	2141	2145	rBV	1878962	3080766	4.49%	1.663%
45	16.010	2145	2148	2155	rVV2	1245892	2650683	3.86%	1.431%
46	16.086	2155	2161	2168	rVB4	1584371	2947935	4.29%	1.592%
47	16.997	2309	2316	2323	rVB	1839967	2754256	4.01%	1.487%
48	17.355	2370	2377	2388	rBV	2294442	3578511	5.21%	1.932%
49	17.555	2404	2411	2417	rBV	2665408	3991110	5.81%	2.155%
50	17.737	2435	2442	2445	rBV	1207624	1868757	2.72%	1.009%
51	17.778	2445	2449	2454	rVV	945951	1419163	2.07%	0.766%
52	17.837	2454	2459	2462	rVV	2233520	3517568	5.12%	1.899%
53	17.872	2462	2465	2472	rVB	1441708	1984138	2.89%	1.071%
54	18.148	2504	2512	2518	rBV	3146129	4706545	6.85%	2.541%
55	18.372	2544	2550	2554	rBV2	135609	196841	0.29%	0.106%
56	19.141	2677	2681	2685	rVB2	97078	124924	0.18%	0.067%
57	19.770	2775	2788	2794	rVV	1930257	2773329	4.04%	1.497%
58	20.105	2838	2845	2848	rVV	2531141	3950690	5.75%	2.133%
59	20.134	2848	2850	2857	rVB	1917933	2214786	3.23%	1.196%
60	20.446	2898	2903	2910	rBV4	122759	178822	0.26%	0.097%
61	22.050	3167	3176	3184	rBV2	2301214	5488279	7.99%	2.963%
62	22.120	3184	3188	3196	rVB	806127	1373400	2.00%	0.741%
63	23.190	3365	3370	3377	rVB2	68493	129672	0.19%	0.070%
64	24.506	3585	3594	3610	rVB	1097450	2979231	4.34%	1.608%
65	25.428	3739	3751	3758	rBV3	1071150	3460100	5.04%	1.868%
66	25.499	3758	3763	3778	rVV2	555932	1602680	2.33%	0.865%
67	25.681	3778	3794	3804	rVB2	524490	1802719	2.63%	0.973%
68	29.870	4489	4507	4515	rBV3	742974	3788761	5.52%	2.046%
69	31.186	4713	4731	4749	rBV4	286620	1647289	2.40%	0.889%

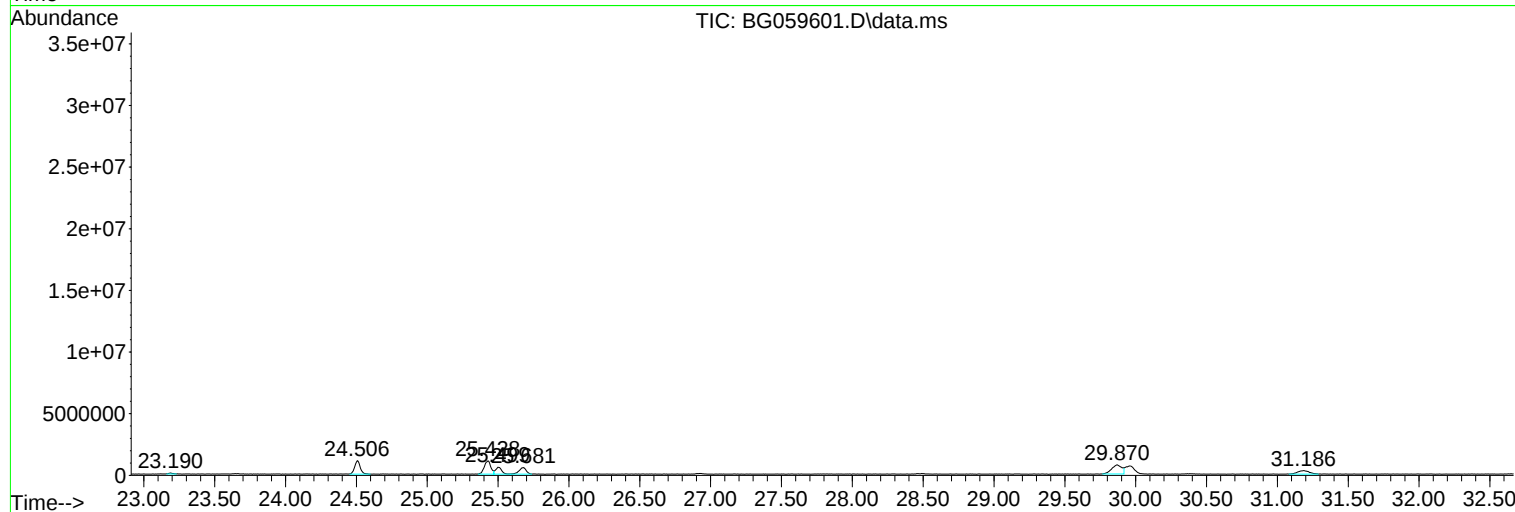
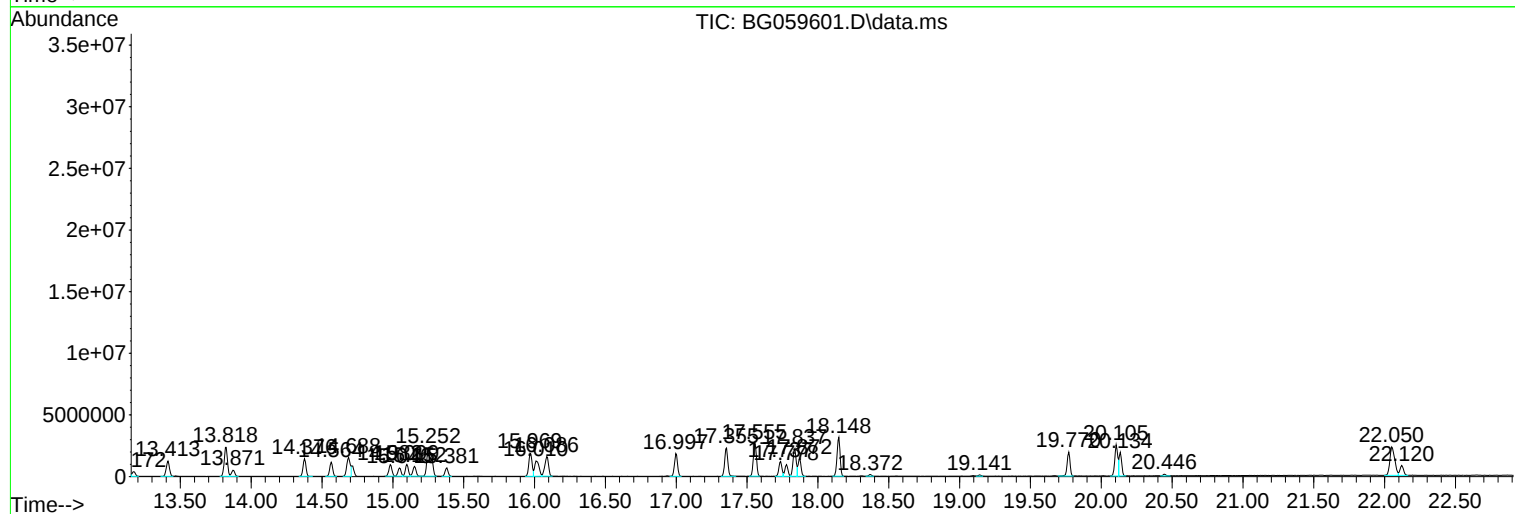
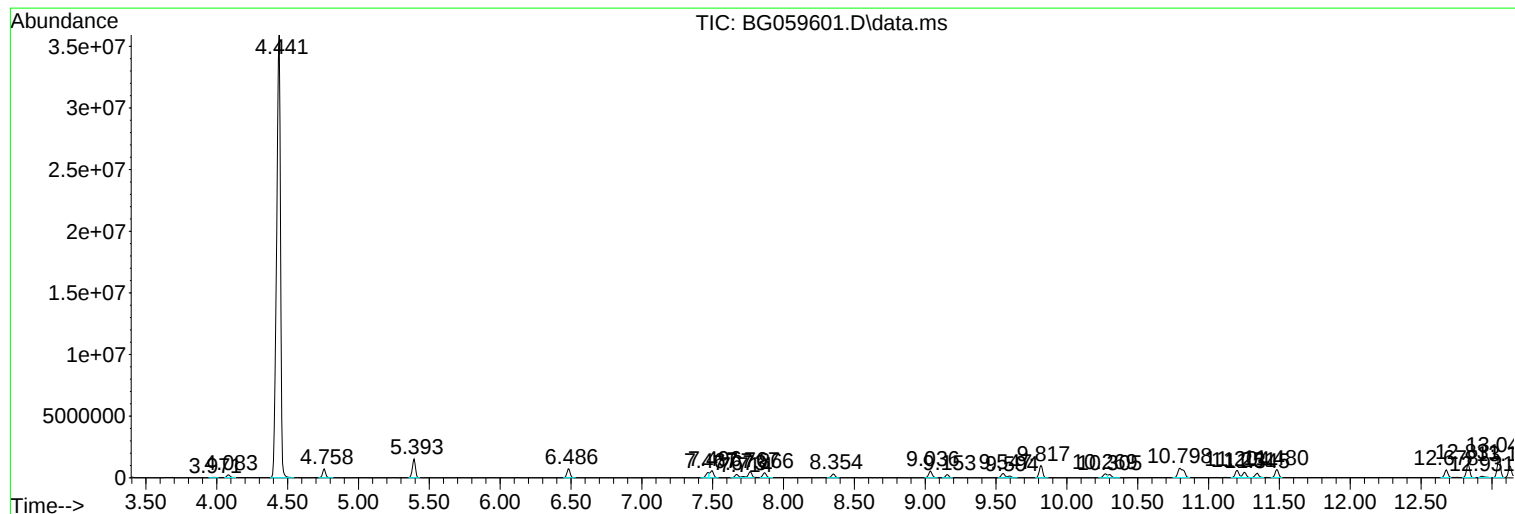
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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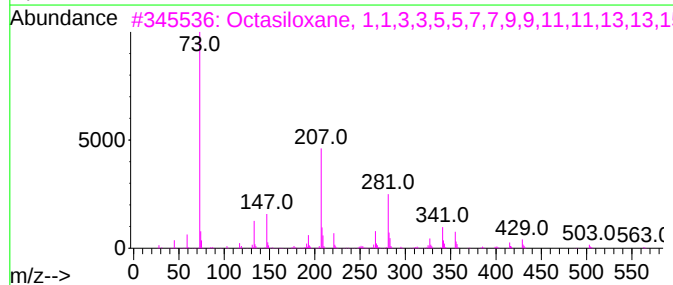
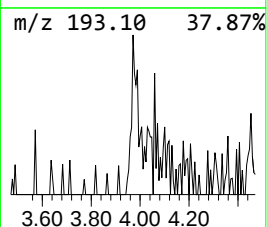
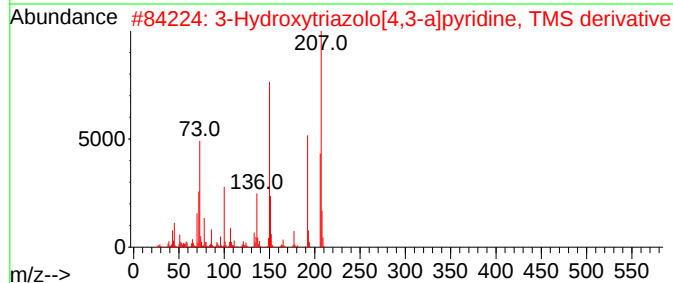
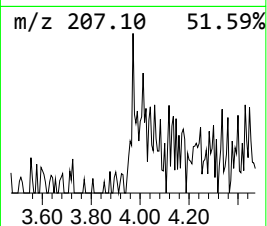
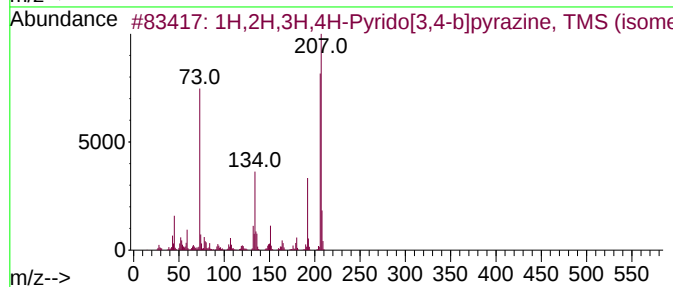
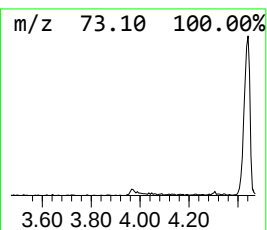
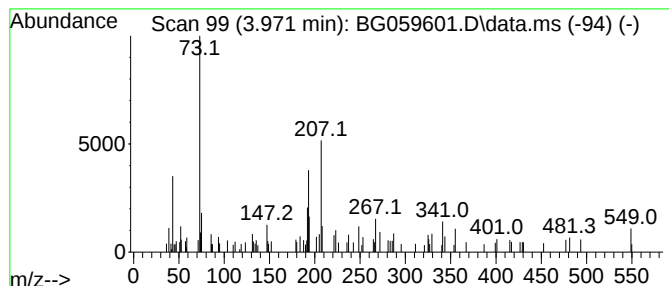
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.971	3.00 ng/ul	77549	1,4-Dichlorobenzene-d4	8.348

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H,2H,3H,4H-Pyrido[3,4-b]pyrazin...	207	C10H17N3Si	1000468-11-7	43
2			3-Hydroxytriazolo[4,3-a]pyridine...	207	C9H13N3OSi	1000476-39-3	43
3			Octasiloxane, 1,1,3,3,5,5,7,7,9,...	578	C16H50O7Si8	019095-24-0	35
4			Dodecahydropyrido[1,2-b]isoquino...	207	C13H21NO	108873-36-5	35
5			Indole-2-one, 2,3-dihydro-N-hydr...	207	C11H13NO3	1010129-52-1	30



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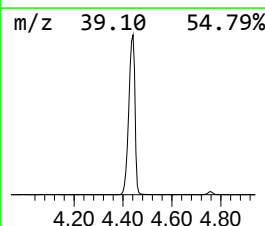
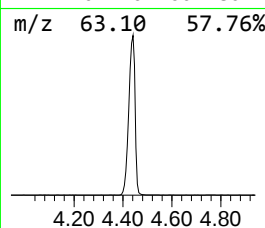
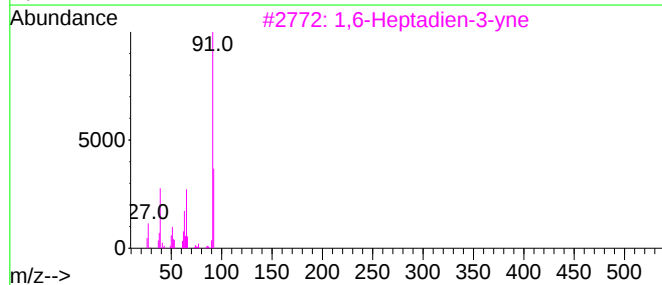
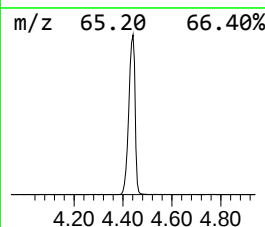
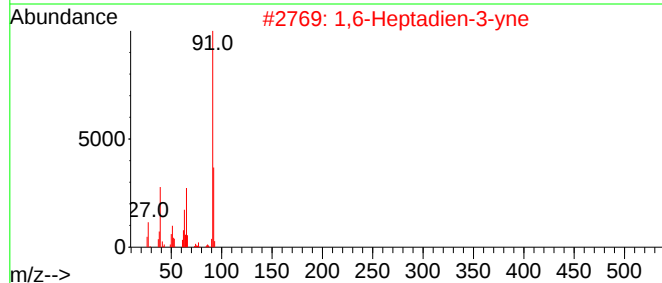
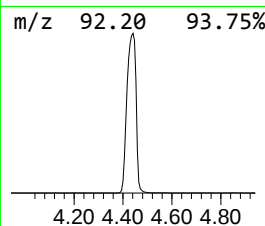
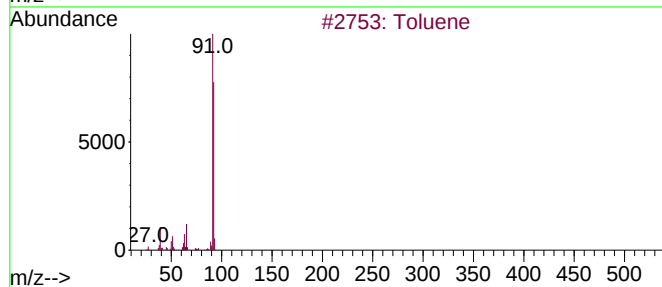
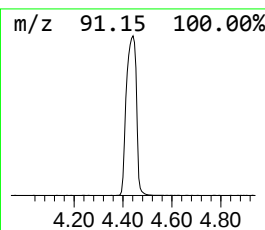
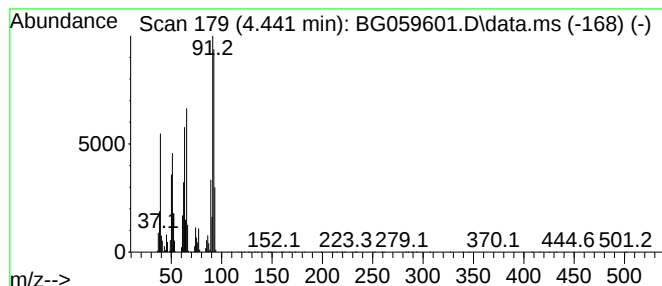
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Toluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.441	2656.37 ng/ul	68674600	1,4-Dichlorobenzene-d4	8.348

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	74
2			1,6-Heptadien-3-yne	92	C7H8	005150-80-1	74
3			1,6-Heptadien-3-yne	92	C7H8	005150-80-1	72
4			Toluene	92	C7H8	000108-88-3	72
5			1,5-Heptadien-3-yne	92	C7H8	003511-27-1	68



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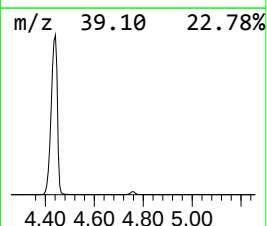
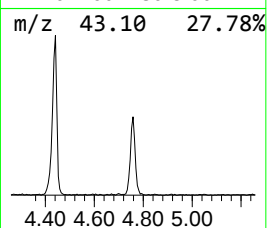
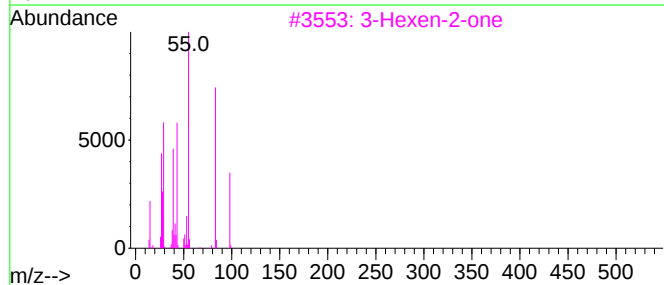
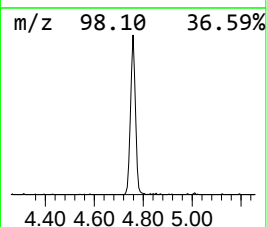
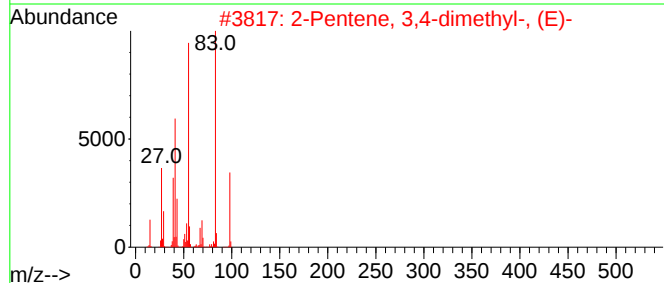
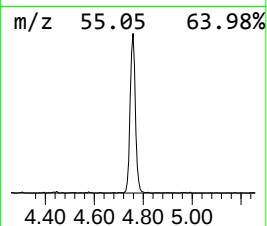
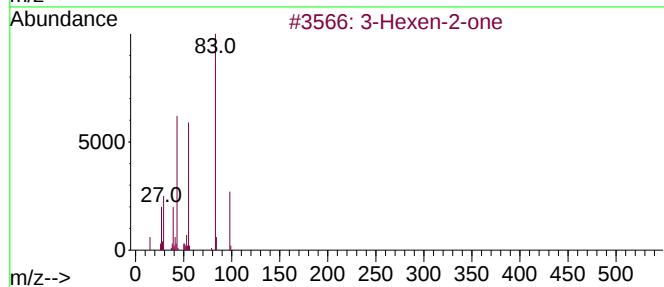
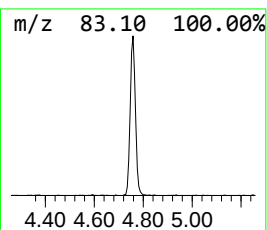
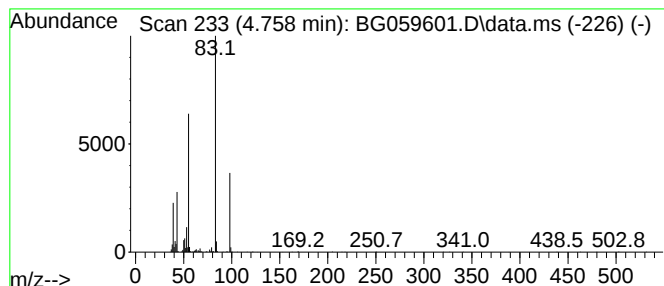
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 3-Hexen-2-one Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.758	40.33 ng/ul	1042560	1,4-Dichlorobenzene-d4	8.348

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Hexen-2-one	98	C6H10O	000763-93-9	90
2			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
3			3-Hexen-2-one	98	C6H10O	000763-93-9	83
4			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	80
5			Furan, 2-methoxy-	98	C5H6O2	025414-22-6	80



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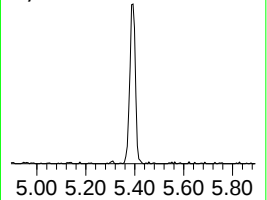
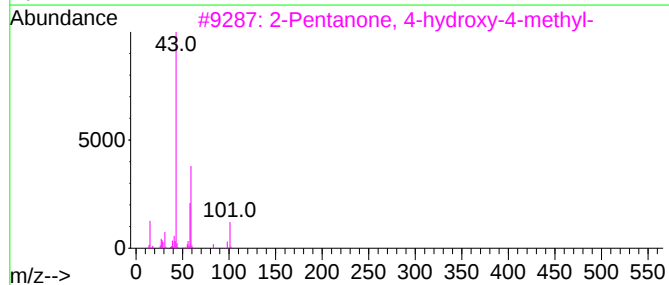
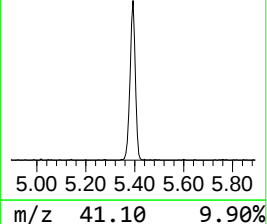
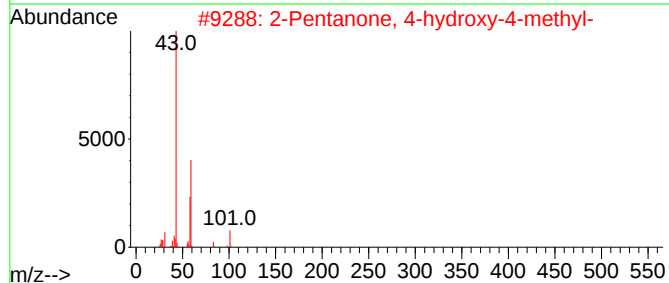
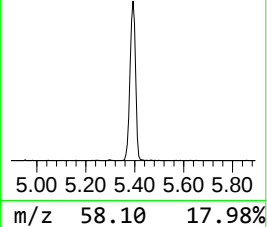
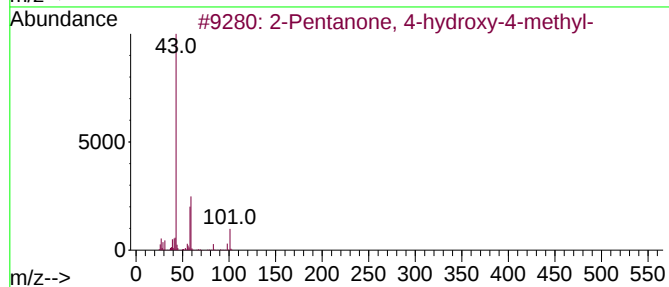
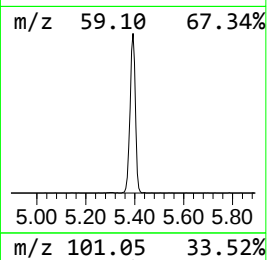
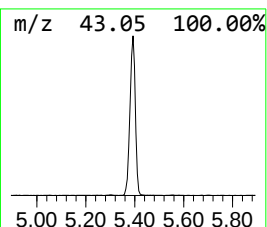
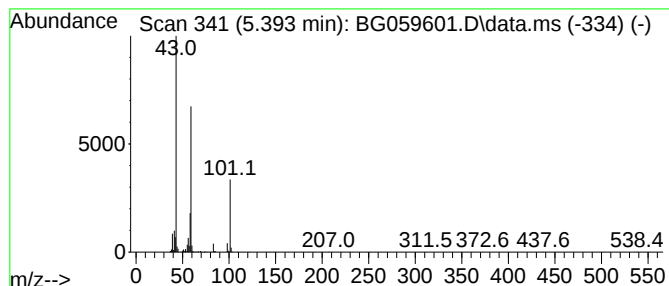
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.393	87.28 ng/ul	2256410	1,4-Dichlorobenzene-d4	8.348

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	50
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

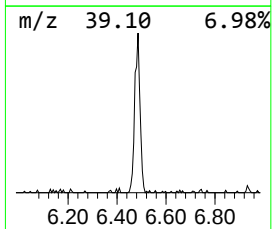
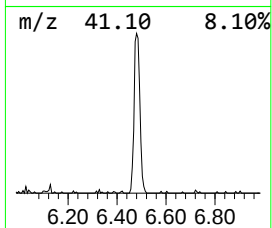
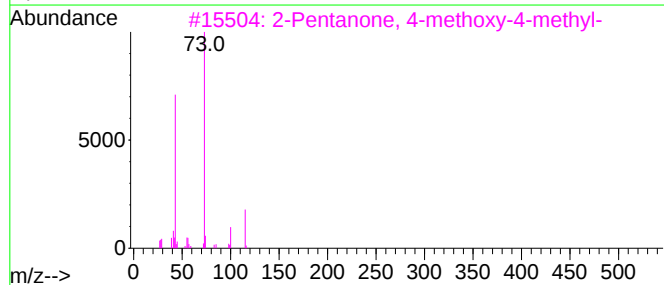
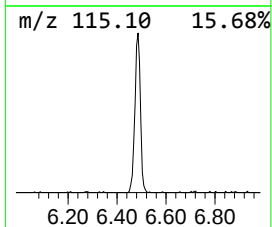
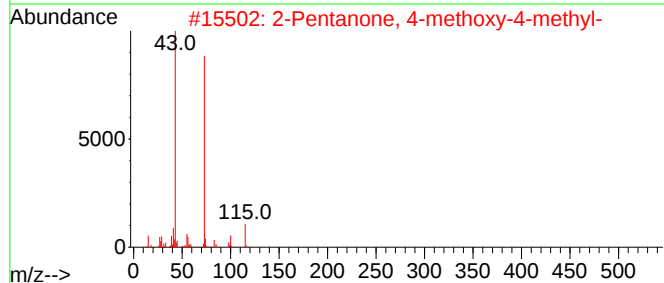
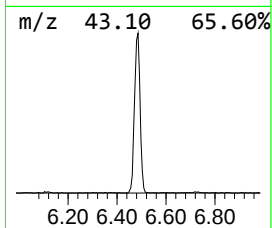
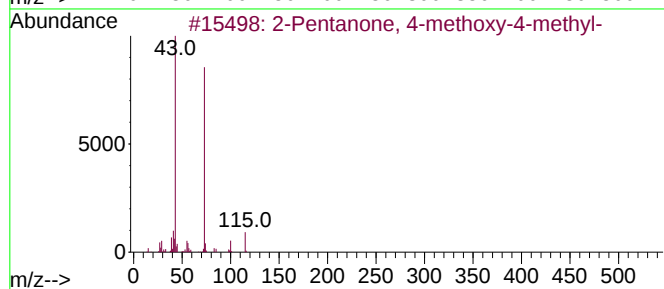
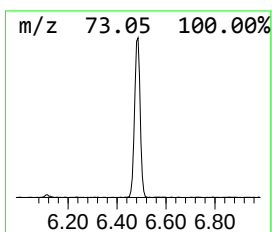
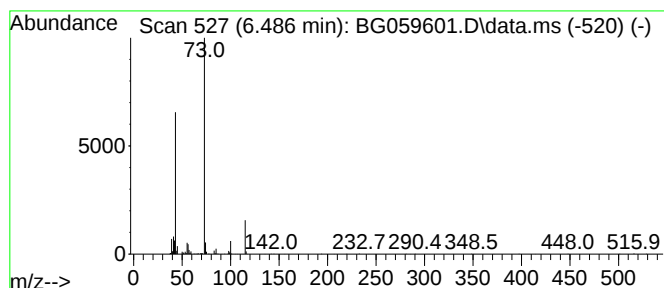
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Pentanone, 4-methoxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.486	44.27 ng/ul	1144500	1,4-Dichlorobenzene-d4			8.348
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-		130	C7H14O2	000107-70-0	83
2	2-Pentanone, 4-methoxy-4-methyl-		130	C7H14O2	000107-70-0	64
3	2-Pentanone, 4-methoxy-4-methyl-		130	C7H14O2	000107-70-0	42
4	Pentane, 3-methoxy-		102	C6H14O	036839-67-5	25
5	Acetamide, N-butyl-		115	C6H13NO	001119-49-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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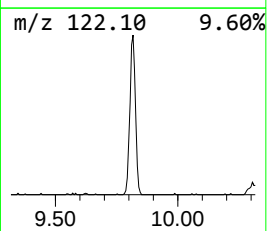
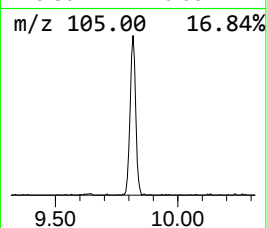
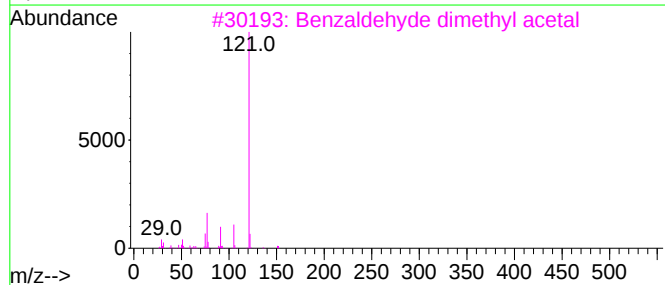
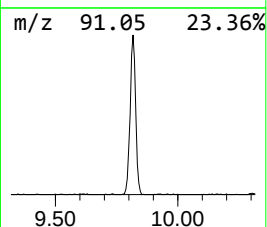
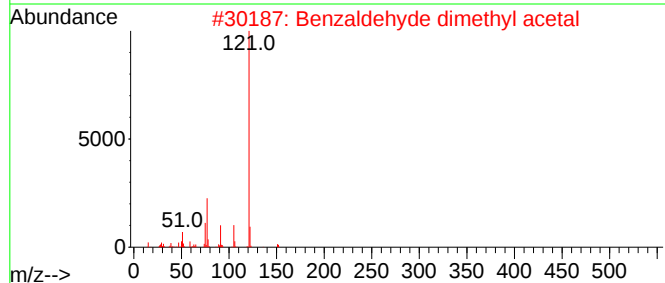
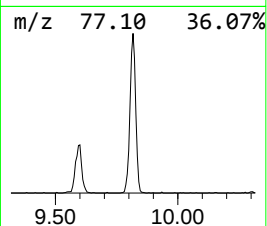
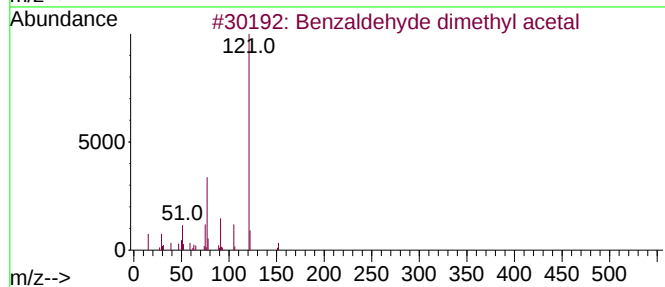
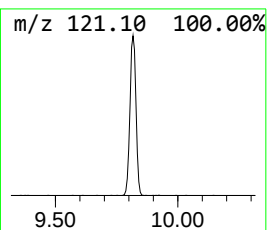
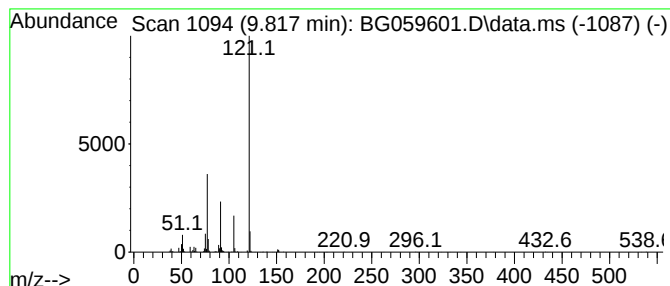
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzaldehyde dimethyl acetal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.817	29.68 ng/ul	1598930	Naphthalene-d8	11.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
2			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
3			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	86
4			Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	83
5			Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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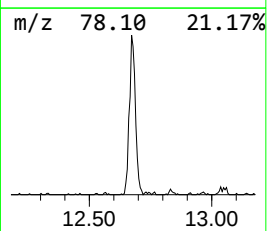
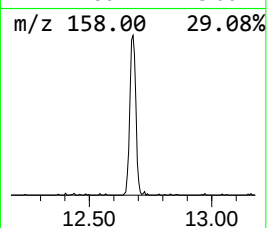
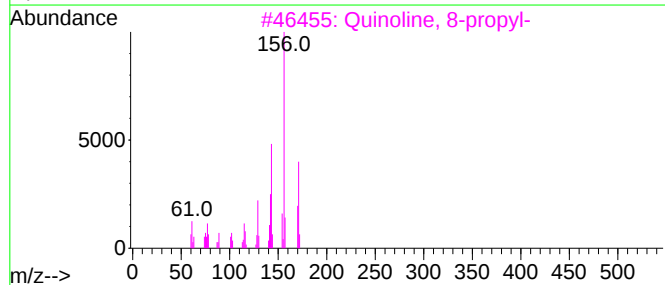
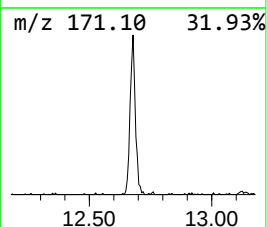
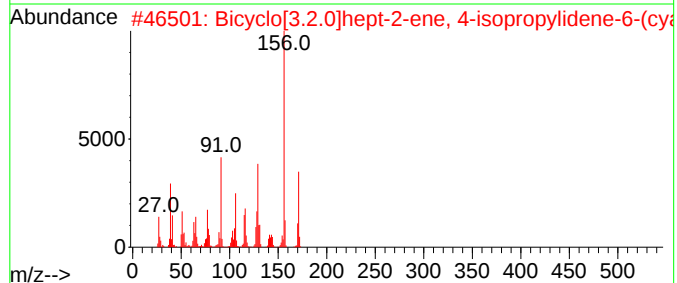
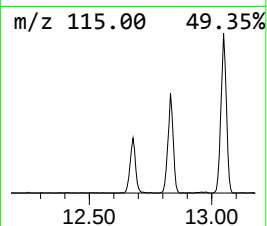
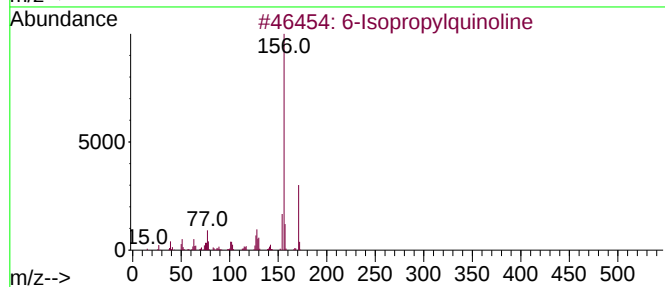
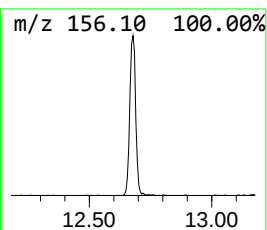
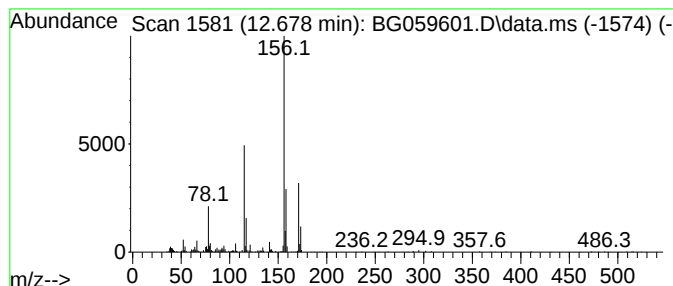
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.678	18.57 ng/ul	1000320	Naphthalene-d8	11.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline		171	C12H13N	000135-79-5	43	
2	Bicyclo[3.2.0]hept-2-ene, 4-isop...		171	C12H13N	1000217-15-6	32	
3	Quinoline, 8-propyl-		171	C12H13N	007661-53-2	32	
4	Cyclopentanone, O-(trimethylsilyl-...		171	C8H17NOSi	026208-87-7	32	
5	Benzene, 2,5-cyclohexadien-1-yl-		156	C12H12	004794-05-2	25	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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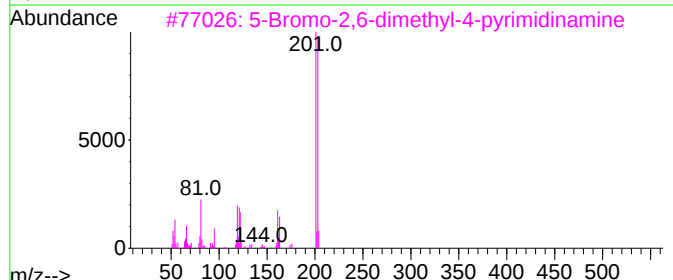
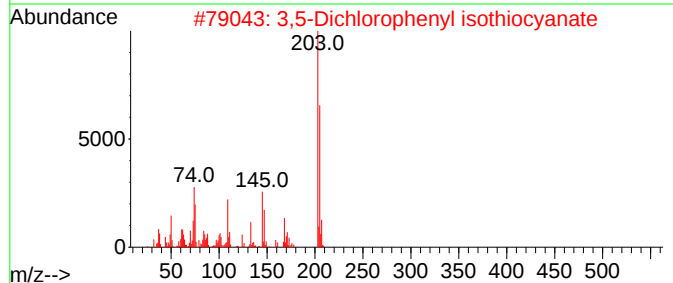
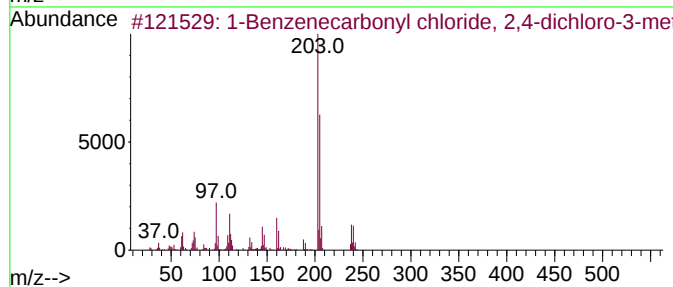
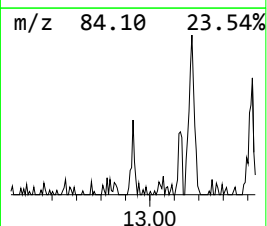
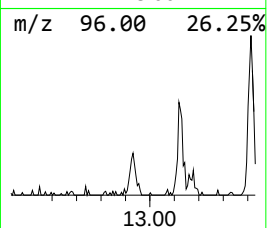
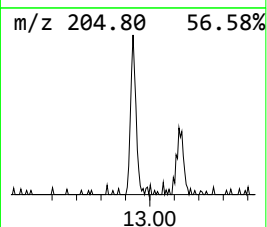
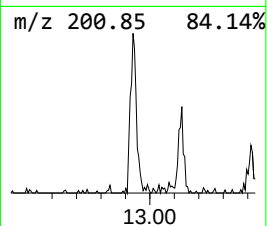
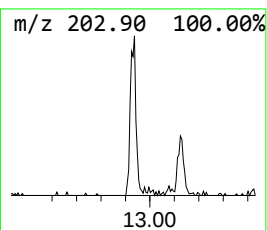
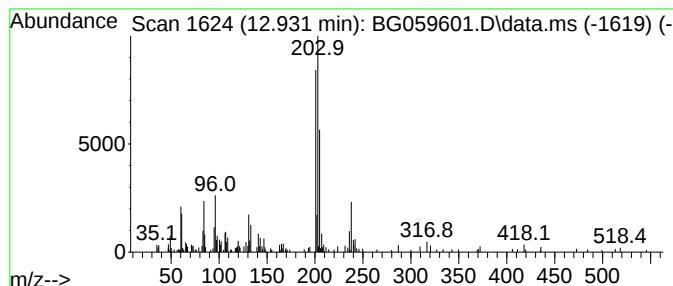
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.931	4.68 ng/ul	252316	Naphthalene-d8	11.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Benzenecarbonyl chloride, 2,4-...	238	C8H5Cl3O2	1000196-85-1	35
2			3,5-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-93-8	30
3			5-Bromo-2,6-dimethyl-4-pyrimidin...	201	C6H8BrN3	007752-80-9	30
4			Imidazo[4,5-b]pyridin-2(3H)-one,...	203	C6H3Cl2N3O	1000269-62-7	27
5			Methyl 3,5-dichloro-4-methoxyben...	234	C9H8Cl2O3	024295-27-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
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Misc :
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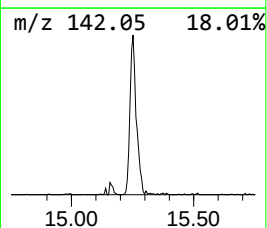
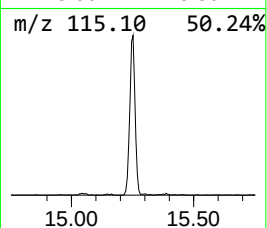
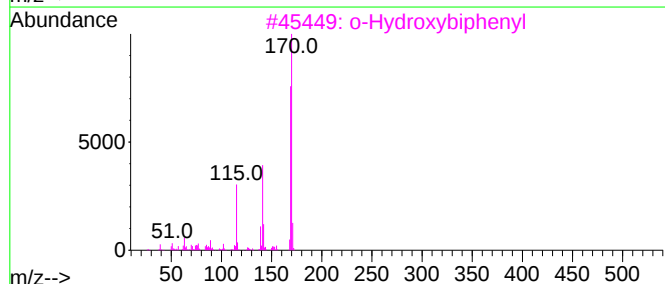
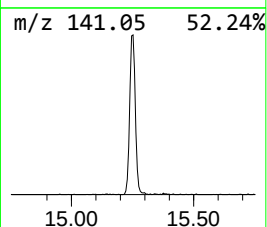
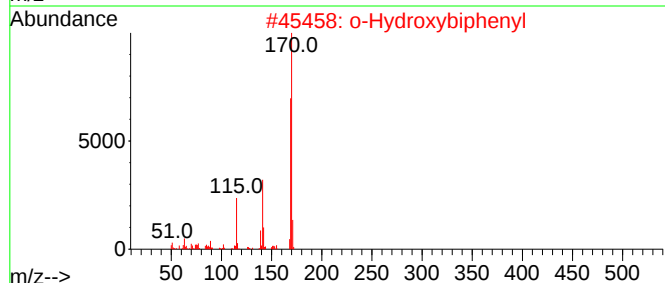
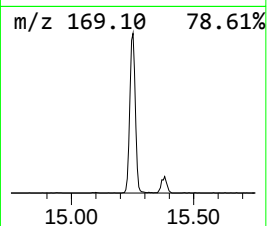
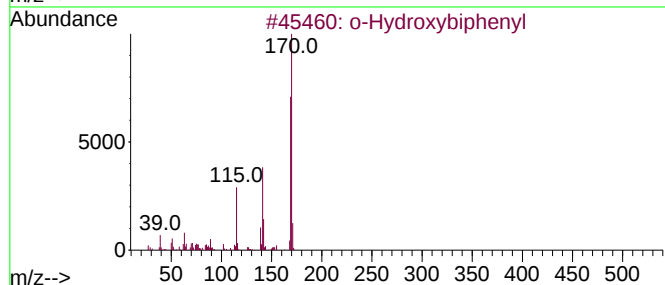
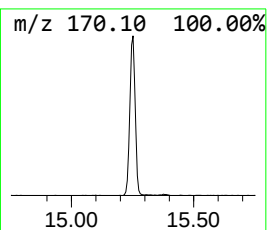
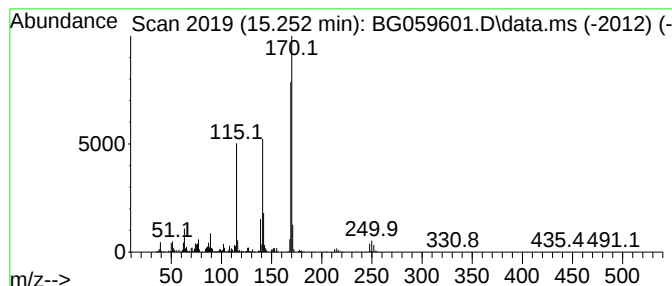
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 o-Hydroxybiphenyl Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.252	76.43 ng/ul	5535710	Acenaphthene-d10	14.982

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
2			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
3			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
4			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	91
5			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
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Sample : 04996-03
Misc :
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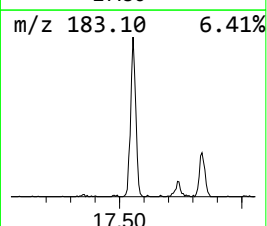
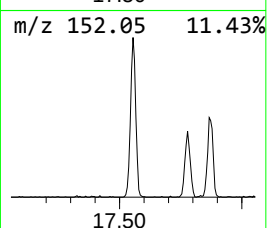
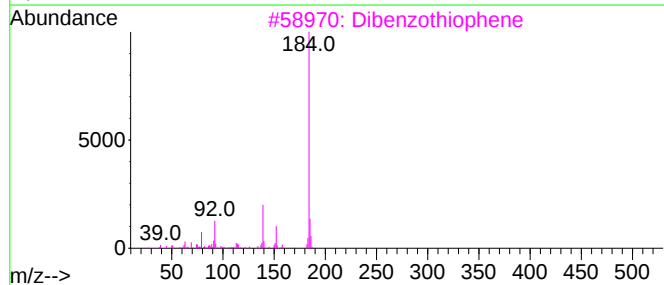
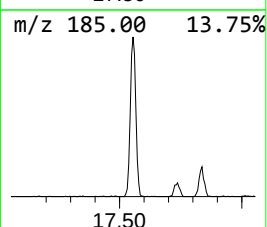
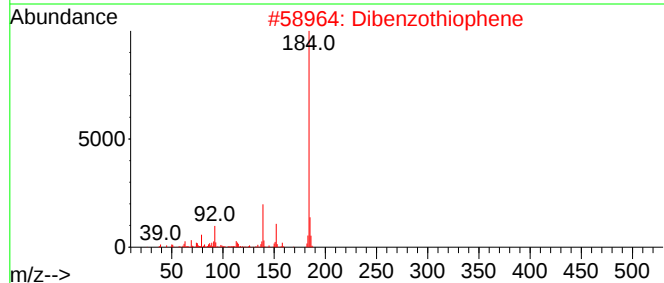
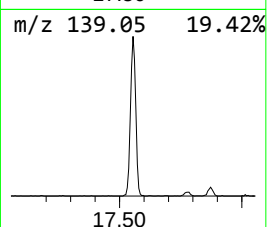
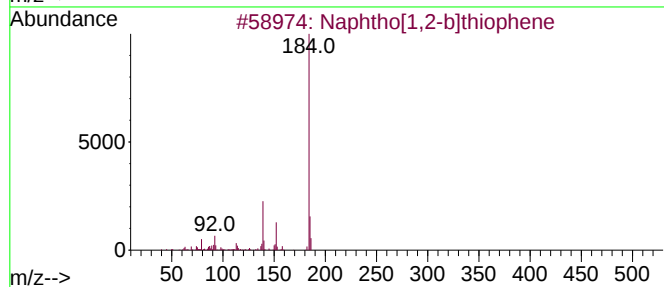
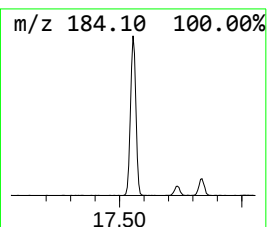
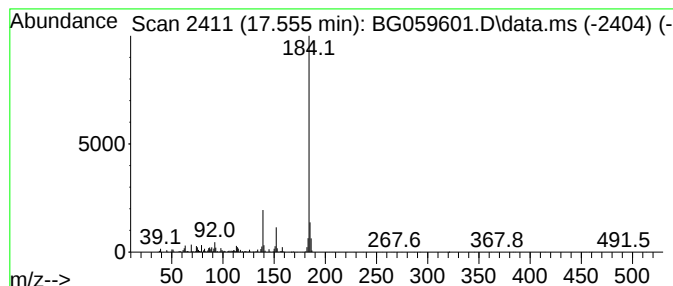
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Naphtho[1,2-b]thiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.555	42.71 ng/ul	3991110	Phenanthrene-d10	17.737

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	97
2			Dibenzothiophene	184	C12H8S	000132-65-0	96
3			Dibenzothiophene	184	C12H8S	000132-65-0	96
4			Dibenzothiophene	184	C12H8S	000132-65-0	96
5			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
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Misc :
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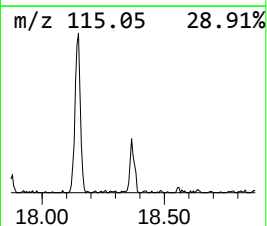
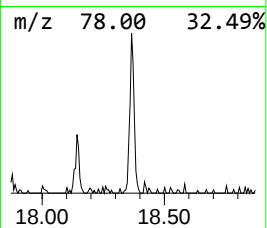
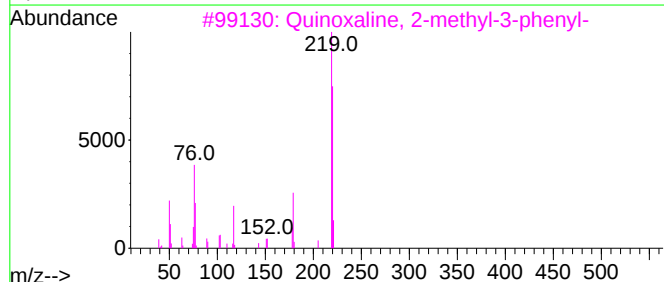
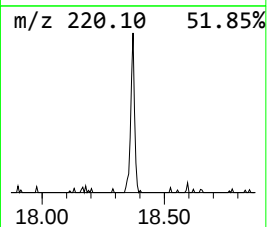
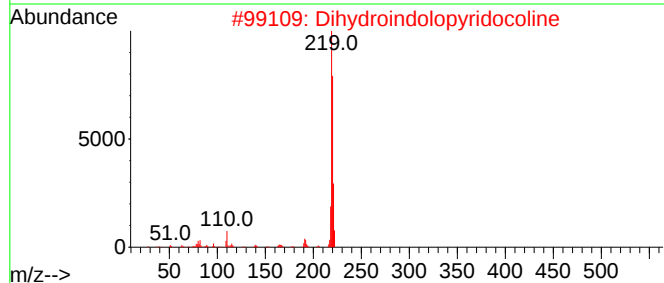
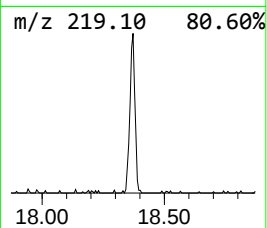
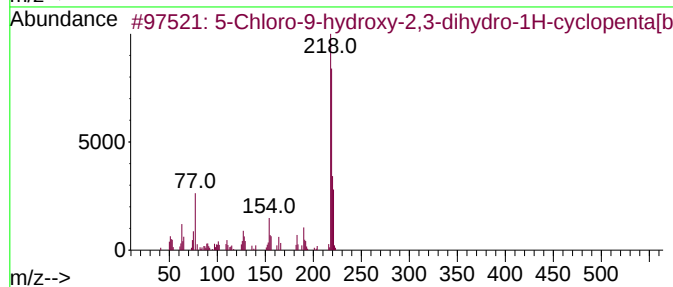
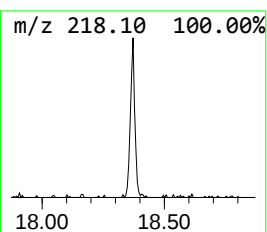
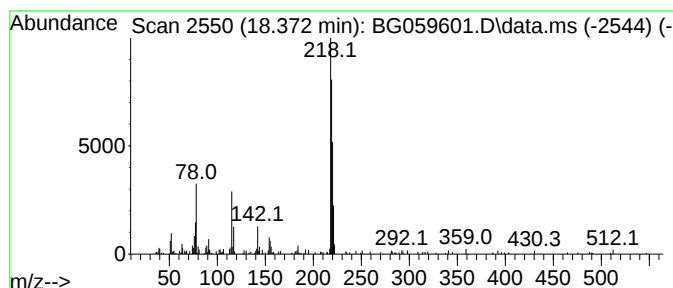
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 5-Chloro-9-hydroxy-2,3-dihy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.372	2.11 ng/ul	196841	Phenanthrene-d10	17.737

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Chloro-9-hydroxy-2,3-dihydro-1...	219	C12H10ClNO	1000254-73-3	72
2			Dihydroindolopyridocoline	220	C15H12N2	117998-81-9	49
3			Quinoxaline, 2-methyl-3-phenyl-	220	C15H12N2	010130-23-1	43
4			N-Phenyl-4-quinolinamine	220	C15H12N2	030696-07-2	43
5			1H-Benzimidazole, 1-(2-phenyleth...	220	C15H12N2	051644-24-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110223\
Data File : BG059601.D
Acq On : 2 Nov 2023 13:33
Operator : MA/JU
Sample : 04996-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4A86

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.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
unknown-01	3.971	3.0	ng/u1	77549	1	8.348	517056	20.0
Toluene	4.441	2656.4	ng/u1	68674600	1	8.348	517056	20.0
3-Hexen-2-one	4.758	40.3	ng/u1	1042560	1	8.348	517056	20.0
2-Pentanone, 4-...	5.393	87.3	ng/u1	2256410	1	8.348	517056	20.0
2-Pentanone, 4-...	6.486	44.3	ng/u1	1144500	1	8.348	517056	20.0
Benzaldehyde di...	9.817	29.7	ng/u1	1598930	2	11.204	1077420	20.0
unknown-02	12.678	18.6	ng/u1	1000320	2	11.204	1077420	20.0
unknown-03	12.931	4.7	ng/u1	252316	2	11.204	1077420	20.0
o-Hydroxybiphenyl	15.252	76.4	ng/u1	5535710	3	14.982	1448630	20.0
Naphtho[1,2-b]t...	17.555	42.7	ng/u1	3991110	4	17.737	1868760	20.0
5-Chloro-9-hydr...	18.372	2.1	ng/u1	196841	4	17.737	1868760	20.0