

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
 Data File : BG062127.D
 Acq On : 18 Jul 2024 00:40
 Operator : MA/JU
 Sample : P3217-22
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YDZJ8

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062127.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.403	15	19	21	rBV3	5197	5484	0.80%	0.061%
2	3.444	22	26	29	rVV5	5236	8603	1.26%	0.095%
3	3.473	29	31	35	rVB4	5946	8140	1.19%	0.090%
4	3.720	67	73	79	rVV	16123	24649	3.61%	0.273%
5	3.861	93	97	108	rBV2	36797	63511	9.30%	0.704%
6	4.014	120	123	125	rBV2	2460	2695	0.39%	0.030%
7	4.202	149	155	158	rBV4	8521	13613	1.99%	0.151%
8	4.625	223	227	229	rBV	2058	2739	0.40%	0.030%
9	5.336	347	348	355	rBV	3451	4106	0.60%	0.045%
10	5.453	366	368	374	rVB3	2967	4811	0.70%	0.053%
11	5.589	388	391	394	rVB2	2162	3175	0.46%	0.035%
12	5.624	394	397	400	rVB2	2124	2746	0.40%	0.030%
13	5.659	400	403	404	rBV2	2829	2825	0.41%	0.031%
14	5.953	449	453	456	rVB2	2550	3004	0.44%	0.033%
15	6.041	462	468	473	rBV3	10252	17285	2.53%	0.192%
16	6.094	473	477	481	rVV3	4561	5958	0.87%	0.066%
17	6.435	533	535	542	rVB	2381	3595	0.53%	0.040%
18	6.981	625	628	630	rBV	3013	3409	0.50%	0.038%
19	7.016	632	634	639	rVV	2355	3133	0.46%	0.035%
20	7.122	645	652	657	rBV5	9556	18284	2.68%	0.203%
21	7.181	657	662	663	rVV	10492	14085	2.06%	0.156%
22	7.222	664	669	679	rVB3	43765	93585	13.70%	1.037%
23	7.363	687	693	699	rBV	71627	116350	17.04%	1.289%
24	7.422	699	703	711	rVB2	22718	36059	5.28%	0.400%
25	7.580	719	730	736	rBV	119847	209490	30.68%	2.321%
26	7.680	742	747	751	rBV3	13320	22227	3.25%	0.246%
27	7.739	754	757	760	rBV	2354	2982	0.44%	0.033%
28	7.851	771	776	778	rBV	2413	3569	0.52%	0.040%
29	8.039	799	808	821	rBV	117060	226135	33.11%	2.505%
30	8.139	821	825	826	rVV2	5095	6277	0.92%	0.070%
31	8.156	826	828	833	rVV4	5254	7135	1.04%	0.079%
32	8.397	863	869	875	rVV2	51187	89844	13.16%	0.995%
33	8.497	883	886	888	rVV2	3175	3500	0.51%	0.039%
34	8.767	924	932	938	rBV2	110623	191287	28.01%	2.119%
35	8.832	941	943	944	rBV2	4493	2707	0.40%	0.030%
36	8.996	966	971	973	rBV2	2316	4485	0.66%	0.050%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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37	9.090	983	987	990	rBV4	3488	4090	0.60%	0.045%
38	9.214	1001	1008	1015	rBV	123398	219774	32.18%	2.435%
39	9.284	1015	1020	1024	rVB3	23214	35181	5.15%	0.390%
40	9.472	1044	1052	1058	rBV5	14432	28355	4.15%	0.314%

41	9.602	1073	1074	1077	rBV2	2823	2931	0.43%	0.032%
42	9.672	1081	1086	1087	rBV2	2854	4718	0.69%	0.052%
43	9.725	1091	1095	1096	rBV	4409	4470	0.65%	0.050%
44	9.754	1096	1100	1104	rVV3	9478	15996	2.34%	0.177%
45	9.790	1104	1106	1109	rVB2	4654	4667	0.68%	0.052%

46	9.931	1124	1130	1137	rBV2	120223	219949	32.21%	2.437%
47	10.030	1142	1147	1155	rVB3	28874	46359	6.79%	0.514%
48	10.160	1167	1169	1173	rBV4	2758	3744	0.55%	0.041%
49	10.201	1173	1176	1178	rBV4	2487	3641	0.53%	0.040%
50	10.260	1181	1186	1192	rVB6	41246	66682	9.76%	0.739%

51	10.330	1197	1198	1201	rBV2	3396	4412	0.65%	0.049%
52	10.401	1205	1210	1212	rBV4	11240	14813	2.17%	0.164%
53	10.442	1214	1217	1218	rBV2	4364	4331	0.63%	0.048%
54	10.483	1218	1224	1231	rVB3	176694	322129	47.17%	3.569%
55	10.547	1233	1235	1238	rBV3	5550	5583	0.82%	0.062%

56	10.647	1250	1252	1257	rVB2	4351	5262	0.77%	0.058%
57	10.694	1257	1260	1261	rBV3	4331	3660	0.54%	0.041%
58	10.853	1280	1287	1293	rBV2	154324	287526	42.10%	3.186%
59	10.900	1293	1295	1301	rVB4	20722	34133	5.00%	0.378%
60	10.947	1301	1303	1305	rBV3	3331	3336	0.49%	0.037%

61	10.994	1305	1311	1318	rBV2	80148	156732	22.95%	1.737%
62	11.217	1344	1349	1353	rBV3	6839	11734	1.72%	0.130%
63	11.364	1369	1374	1377	rBV4	7228	11016	1.61%	0.122%
64	11.435	1383	1386	1388	rBV4	7492	6521	0.95%	0.072%
65	11.587	1406	1412	1421	rVB3	23576	50705	7.42%	0.562%

66	11.652	1421	1423	1426	rBV4	2550	3870	0.57%	0.043%
67	11.681	1426	1428	1429	rBV2	3471	2782	0.41%	0.031%
68	11.764	1438	1442	1448	rVB8	9675	18838	2.76%	0.209%
69	11.805	1448	1449	1453	rVB3	5135	3800	0.56%	0.042%
70	11.881	1458	1462	1467	rBV6	8067	13474	1.97%	0.149%

71	11.952	1470	1474	1477	rBV4	8808	12782	1.87%	0.142%
72	12.034	1484	1488	1493	rVB5	5694	10349	1.52%	0.115%
73	12.169	1505	1511	1512	rBV5	16758	24900	3.65%	0.276%
74	12.193	1512	1515	1519	rVB3	34819	45188	6.62%	0.501%
75	12.304	1530	1534	1538	rVB7	5636	7429	1.09%	0.082%

76	12.510	1563	1569	1575	rVB	65536	112781	16.51%	1.250%
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77	12.657	1592	1594	1596	rBV2	4912	2830	0.41%	0.031%
78	12.727	1600	1606	1617	rVV	181552	322488	47.22%	3.573%
79	12.821	1620	1622	1628	rVB7	9591	17645	2.58%	0.195%
80	13.256	1693	1696	1702	rBV7	7930	14181	2.08%	0.157%
81	13.321	1705	1707	1708	rBV2	6144	4298	0.63%	0.048%
82	13.468	1731	1732	1735	rBV3	5983	5052	0.74%	0.056%
83	13.532	1738	1743	1746	rBV7	9245	18100	2.65%	0.201%
84	13.838	1793	1795	1797	rBV3	9660	9940	1.46%	0.110%
85	13.938	1810	1812	1813	rBV2	6117	2793	0.41%	0.031%
86	13.996	1818	1822	1827	rVV4	33214	60572	8.87%	0.671%
87	14.079	1829	1836	1845	rVB	301732	496355	72.68%	5.499%
88	14.372	1880	1886	1898	rVB	321738	519159	76.02%	5.752%
89	14.678	1932	1938	1945	rVB	262783	414440	60.69%	4.592%
90	14.913	1974	1978	1986	rVB4	72807	130495	19.11%	1.446%
91	15.130	2011	2015	2022	rVB3	41430	68710	10.06%	0.761%
92	15.406	2058	2062	2066	rVB	56284	78008	11.42%	0.864%
93	15.665	2101	2106	2113	rVB2	415402	622874	91.21%	6.901%
94	15.788	2123	2127	2133	rVB2	149151	204911	30.01%	2.270%
95	16.611	2264	2267	2271	rBV4	33609	43843	6.42%	0.486%
96	17.422	2400	2405	2411	rVB2	329998	487393	71.37%	5.400%
97	17.522	2417	2422	2429	rVB	438248	637474	93.35%	7.063%
98	19.807	2806	2811	2817	rBV	495366	682911	100.00%	7.566%
99	21.681	3126	3130	3137	rVB	292918	463645	67.89%	5.137%
100	24.684	3636	3641	3653	rVB2	249093	647335	94.79%	7.172%

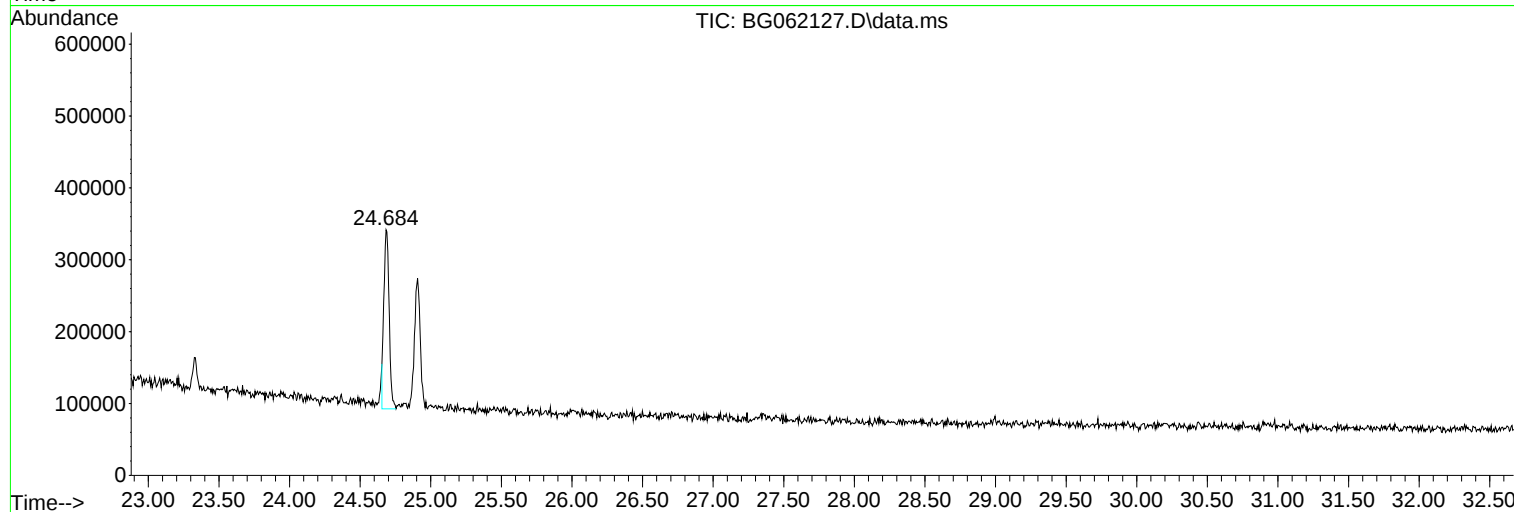
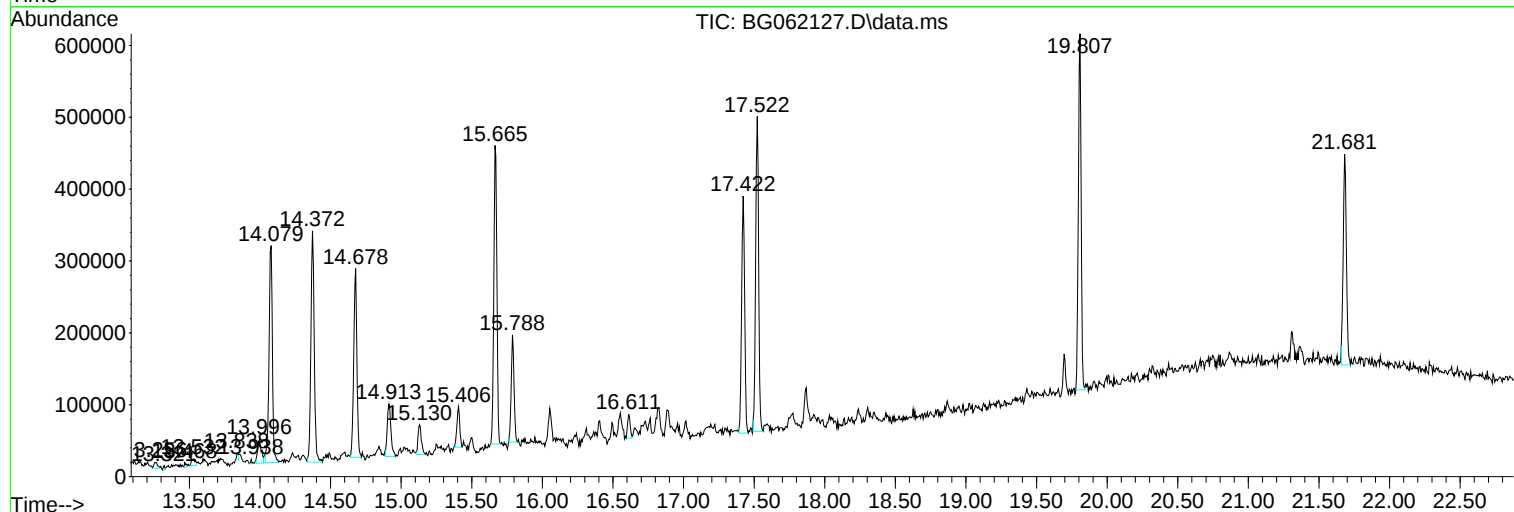
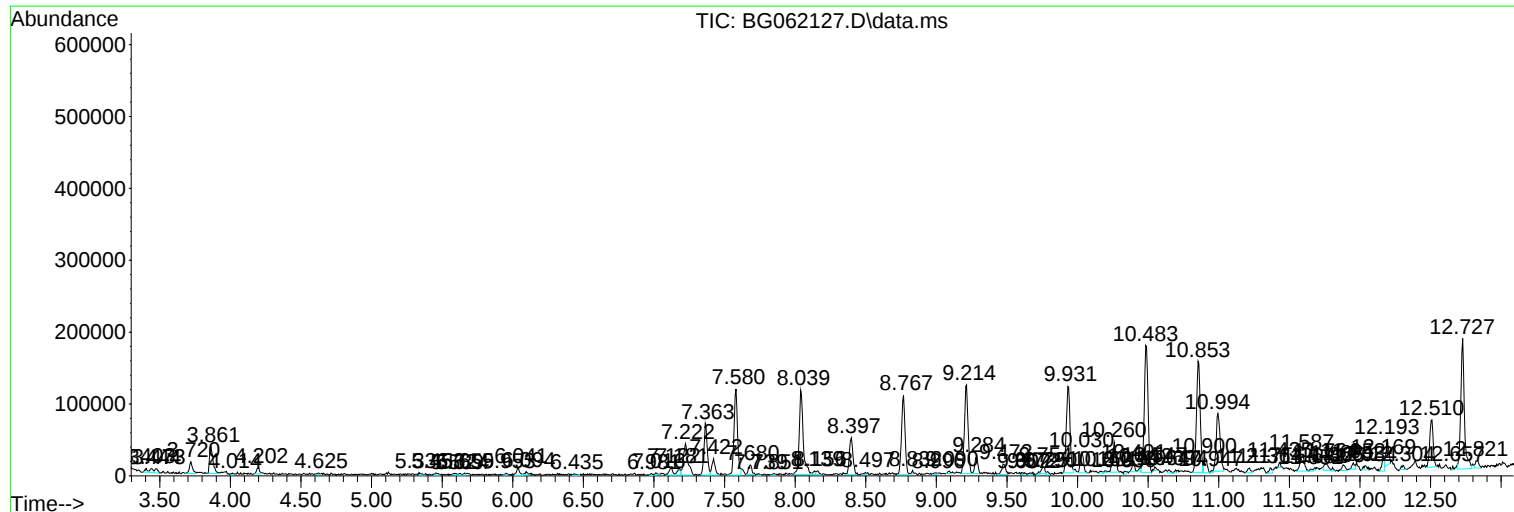
Sum of corrected areas: 9025577

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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



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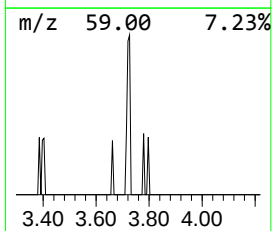
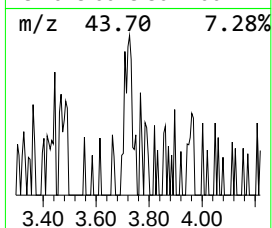
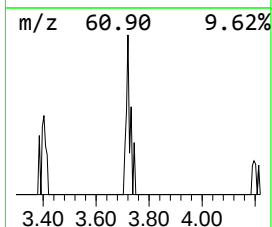
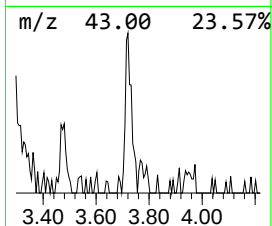
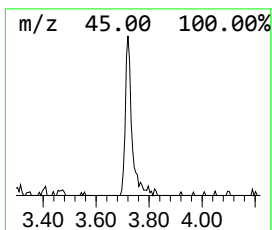
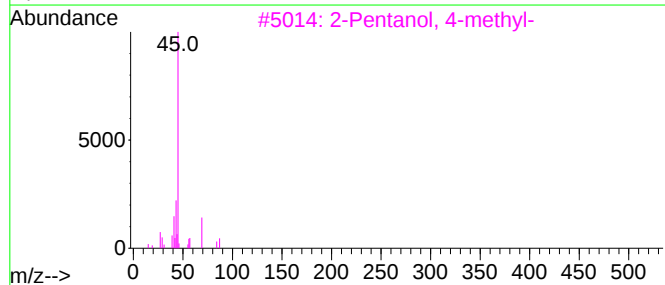
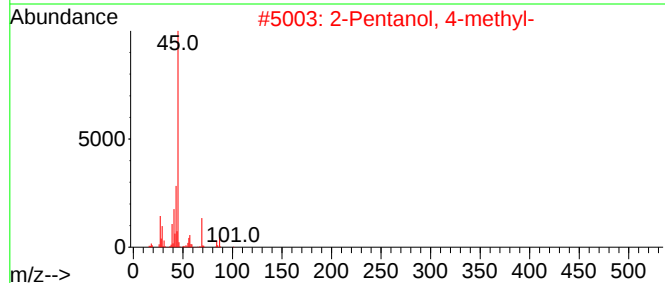
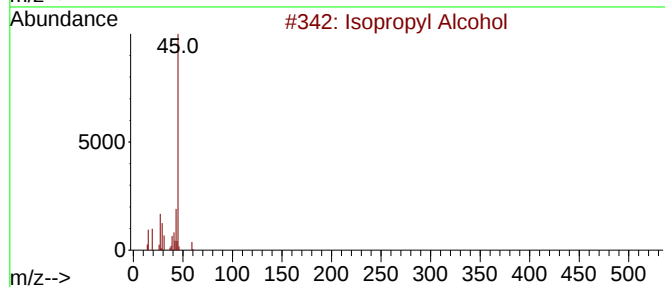
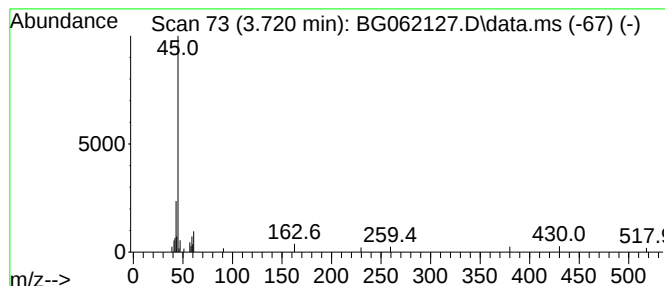
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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.
3.720	2.18 ng/ul	24649	1,4-Dichlorobenzene-d4	8.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	36
2			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	33
3			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	33
4			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	17
5			d-Arabinose, cyclic 1,2-ethanedi...	394	C15H22O8S2	040733-16-2	10



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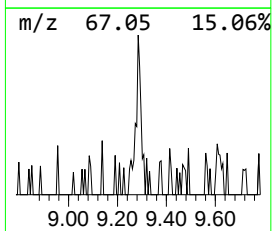
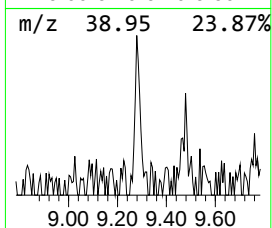
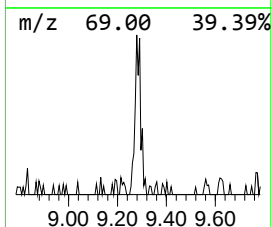
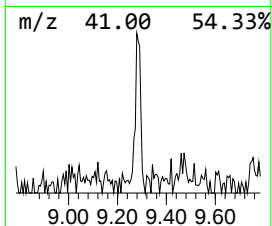
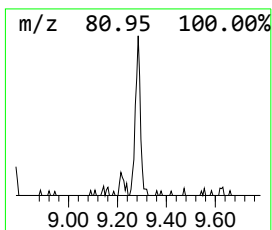
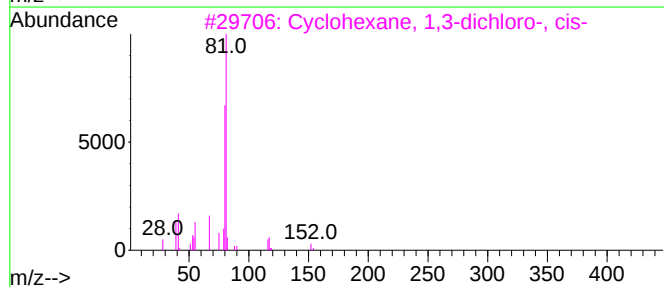
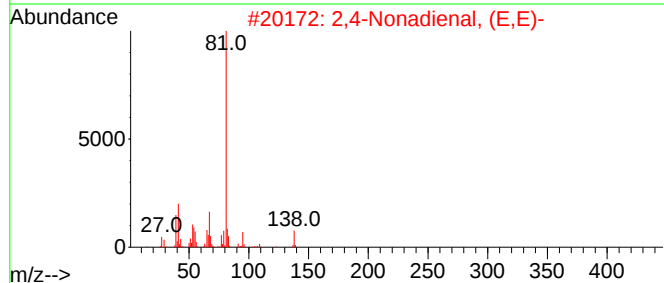
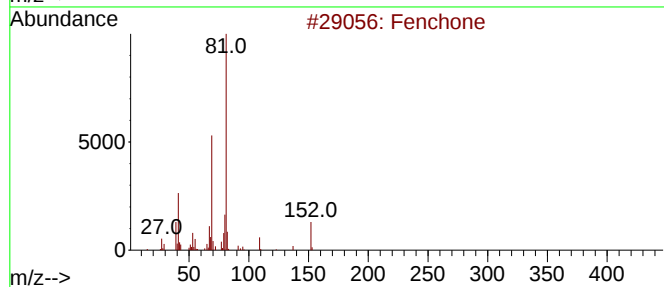
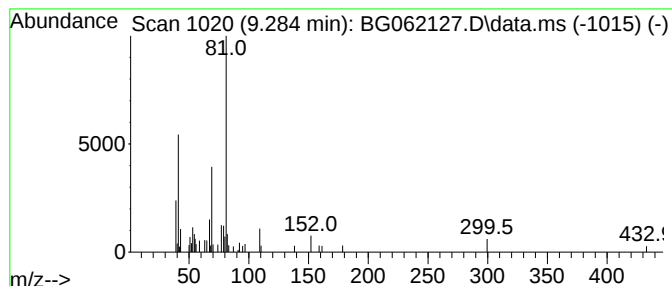
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Fenchone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.284	3.11 ng/ul	35181	1,4-Dichlorobenzene-d4	8.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchone	152	C10H16O	001195-79-5	64
2			2,4-Nonadienal, (E,E)-	138	C9H14O	005910-87-2	53
3			Cyclohexane, 1,3-dichloro-, cis-	152	C6H10Cl2	024955-63-3	53
4			2-Furanacetaldehyde, .alpha.-pro...	152	C9H12O2	031681-26-2	52
5			L-Fenchone	152	C10H16O	007787-20-4	45



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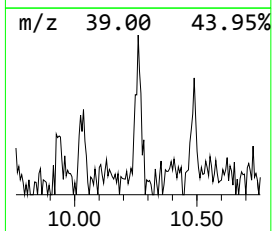
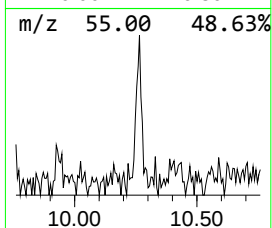
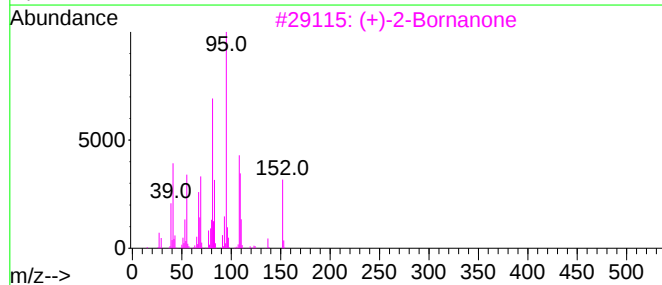
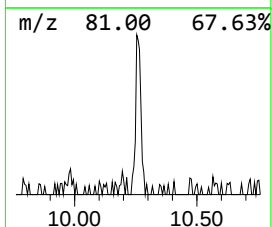
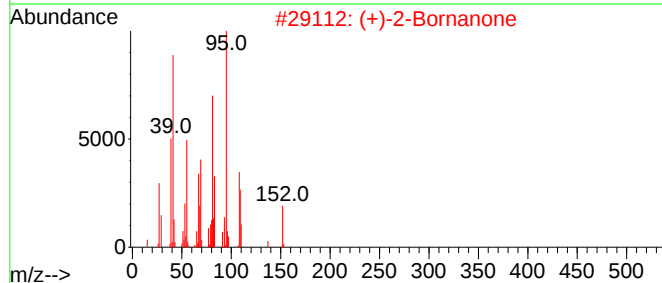
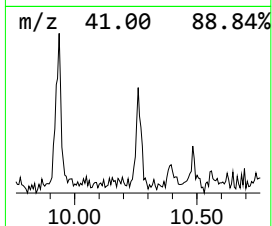
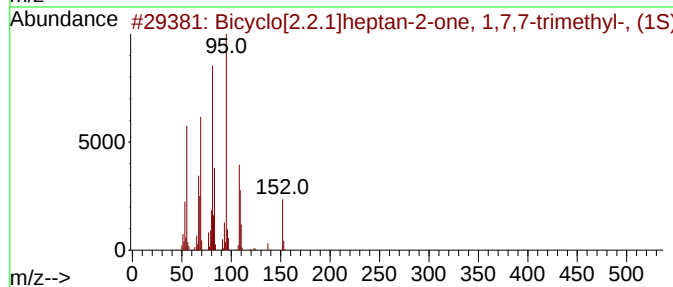
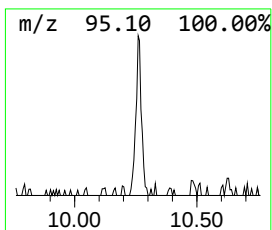
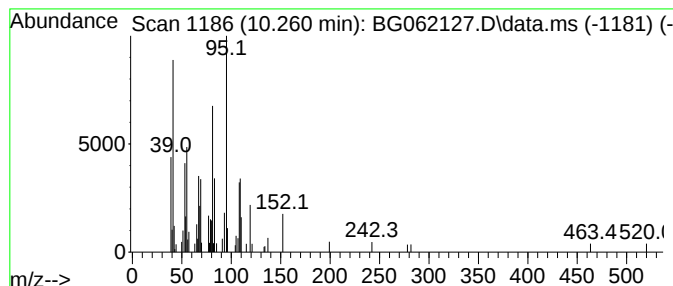
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.260	4.64 ng/ul	66682	Naphthalene-d8	10.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	83
2			(+)-2-Bornanone	152	C10H16O	000464-49-3	80
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	55
4			(+)-2-Bornanone	152	C10H16O	000464-49-3	49
5			3,4-Pentadienal, 2,2-dimethyl-	110	C7H10O	004058-51-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZJ8

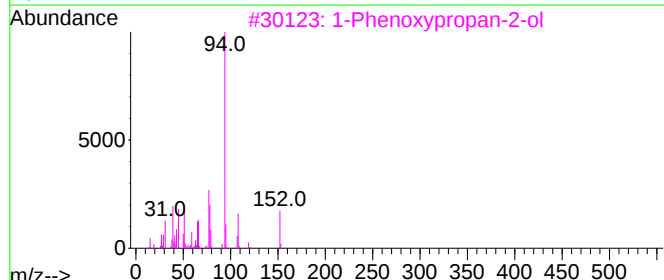
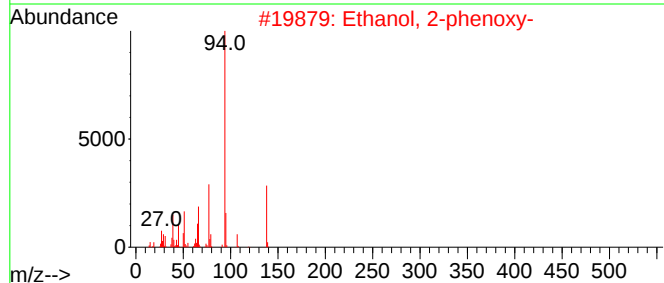
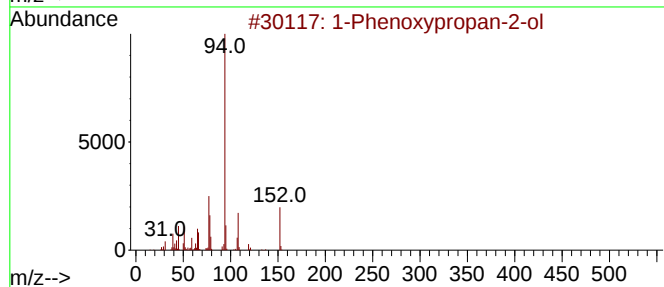
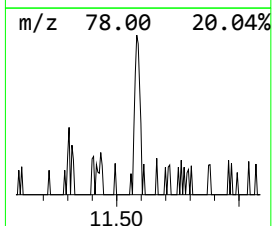
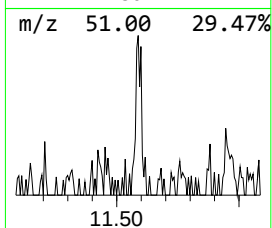
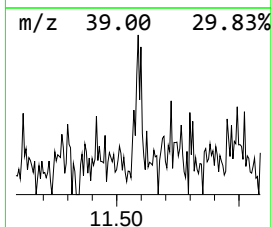
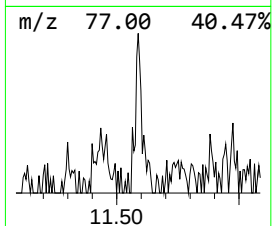
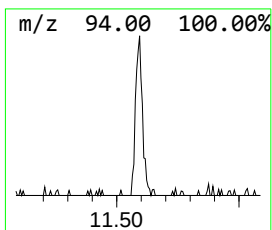
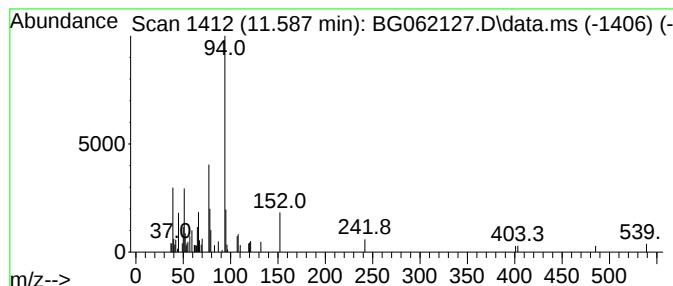
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.587	3.53 ng/ul	50705	Naphthalene-d8	10.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64
2			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64
4			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZJ8

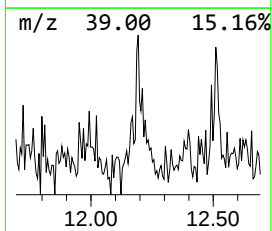
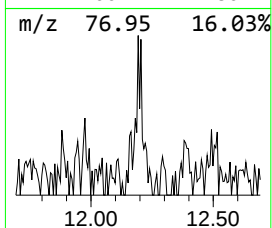
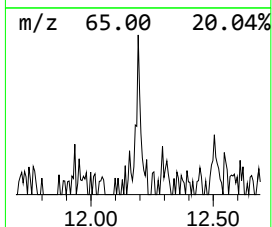
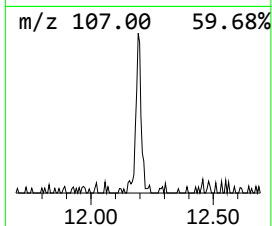
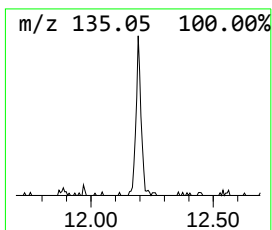
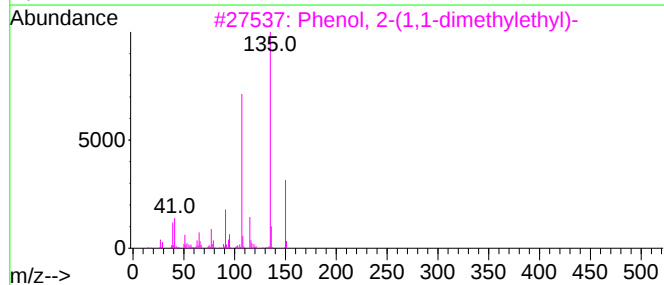
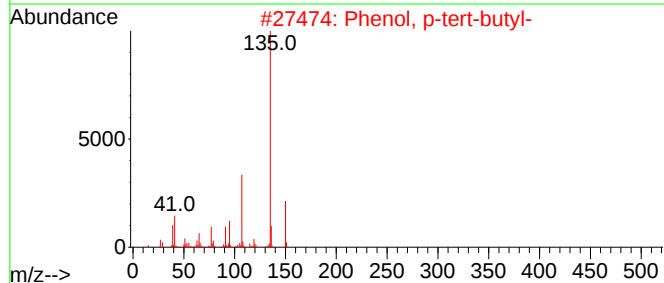
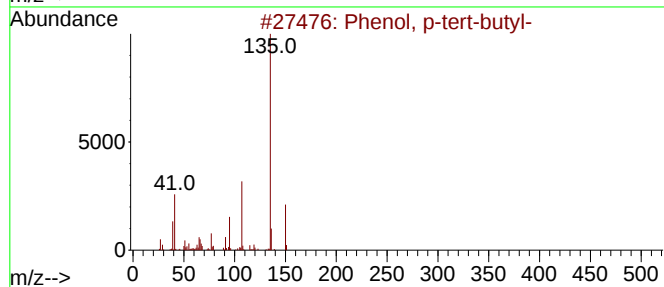
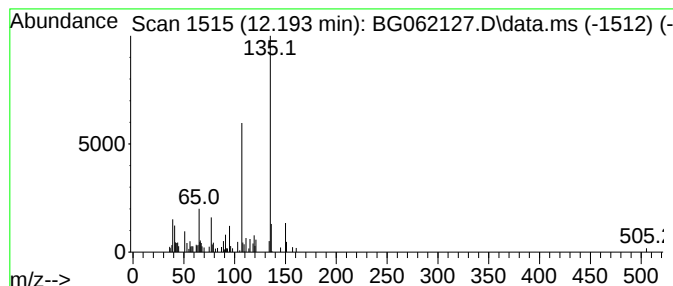
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, p-tert-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.193	3.14 ng/ul	45188	Naphthalene-d8	10.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3			Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	64
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	59
5			4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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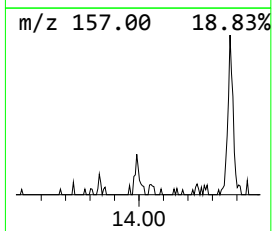
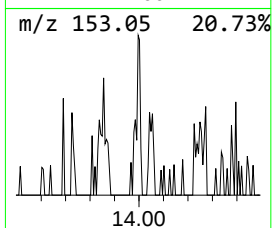
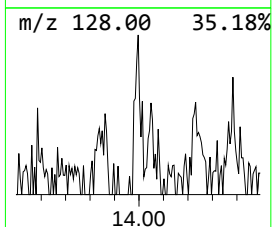
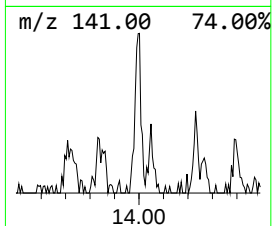
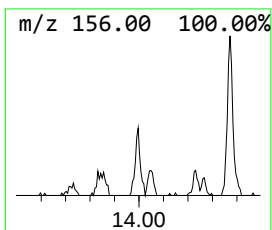
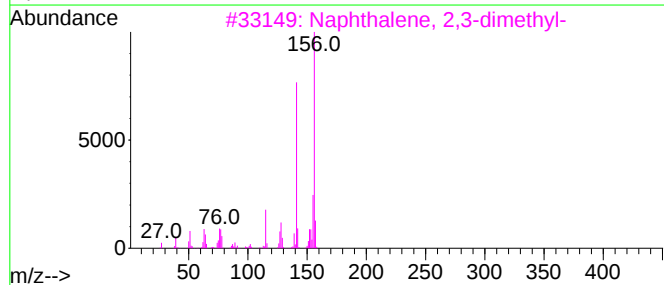
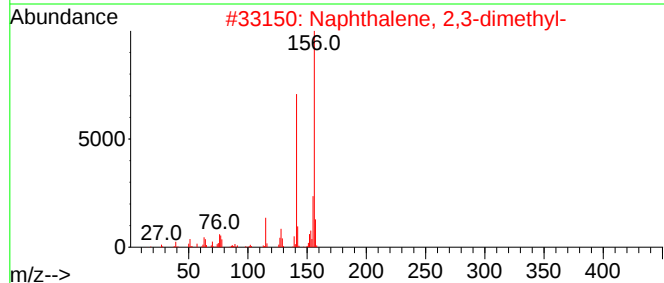
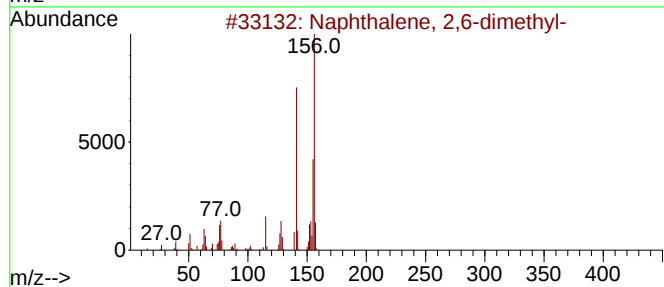
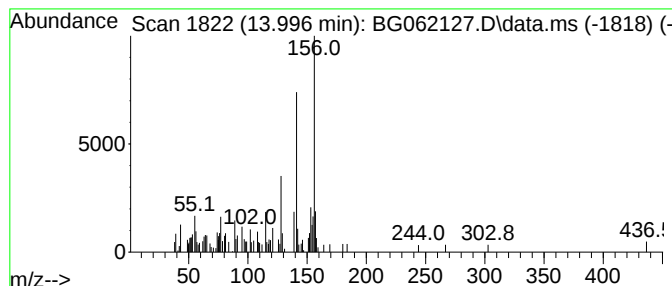
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.996	2.92 ng/ul	60572	Acenaphthene-d10	14.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	81
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	81
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	74
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	72
5			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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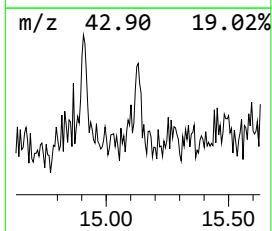
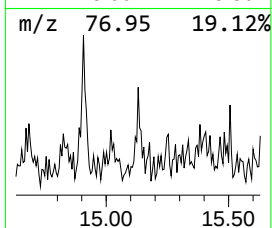
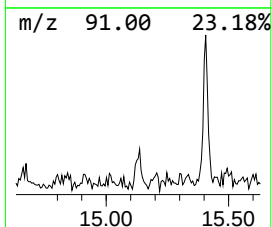
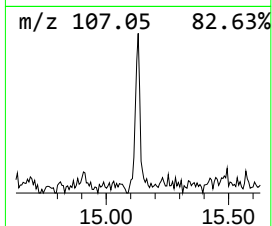
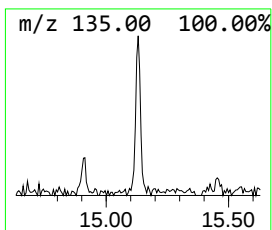
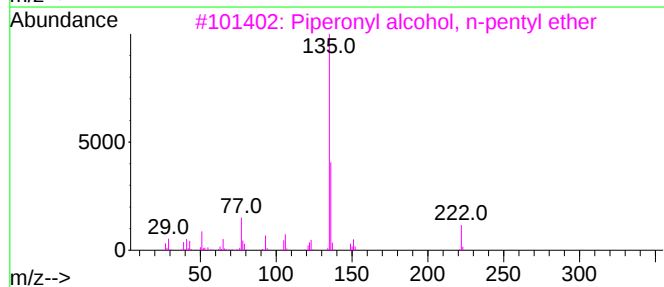
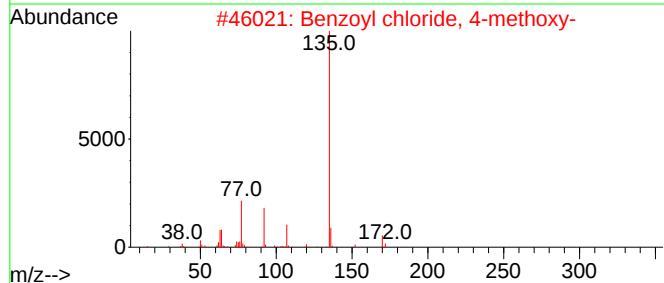
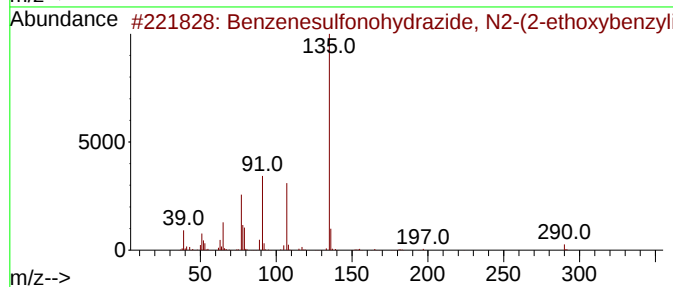
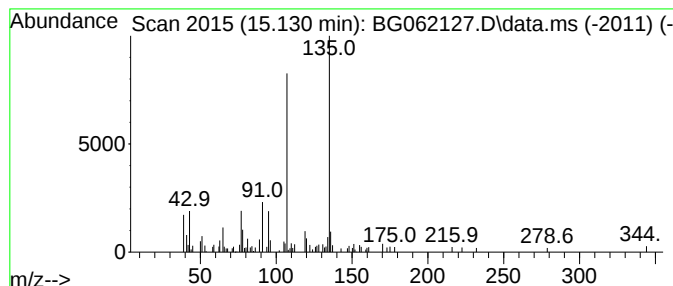
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.130	3.32 ng/ul	68710	Acenaphthene-d10	14.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonohydrazide, N2-(2-e...	318	C16H18N2O3S	065609-76-9	43
2			Benzoyl chloride, 4-methoxy-	170	C8H7ClO2	000100-07-2	43
3			Piperonyl alcohol, n-pentyl ether	222	C13H18O3	1000395-11-6	41
4			4-Methoxy-N'-(2-oxoindolin-3-yl...	295	C16H13N3O3	082973-18-0	38
5			Benzoic acid, 4-methoxy-, methyl...	166	C9H10O3	000121-98-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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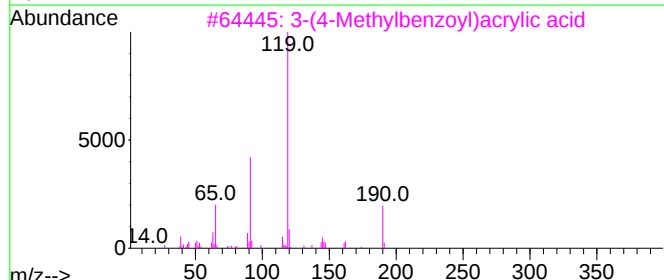
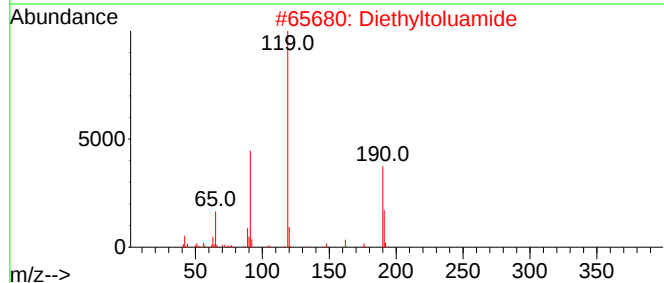
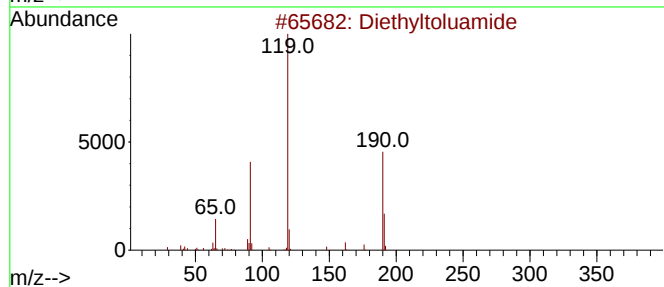
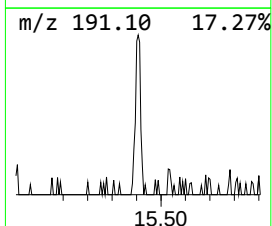
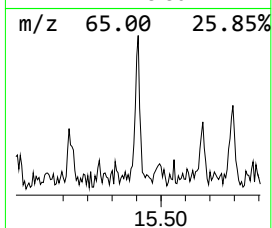
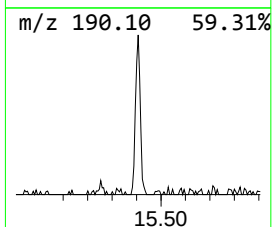
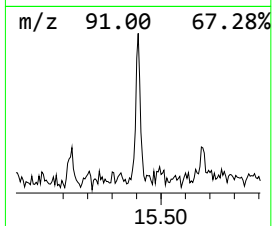
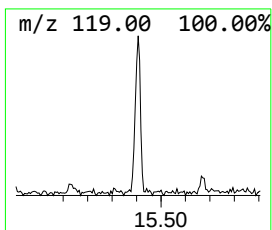
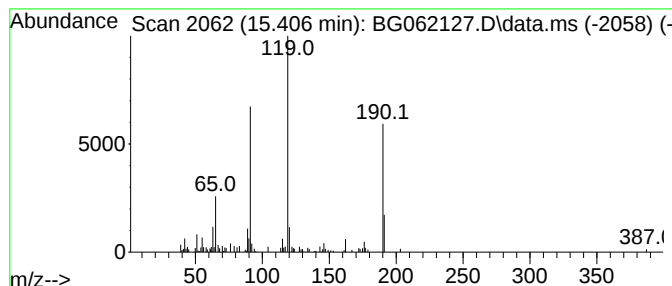
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Diethyltoluamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.406	3.76 ng/ul	78008	Acenaphthene-d10	14.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyltoluamide	191	C12H17NO	000134-62-3	91
2			Diethyltoluamide	191	C12H17NO	000134-62-3	74
3			3-(4-Methylbenzoyl)acrylic acid	190	C11H10O3	020972-36-5	72
4			Diethyltoluamide	191	C12H17NO	000134-62-3	64
5			Diethyltoluamide	191	C12H17NO	000134-62-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071724\
Data File : BG062127.D
Acq On : 18 Jul 2024 00:40
Operator : MA/JU
Sample : P3217-22
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZJ8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Fenchone	9.284	3.1	ng/ul	35181	1	8.045	226135	20.0
Bicyclo[2.2.1]h...	10.260	4.6	ng/ul	66682	2	10.853	287526	20.0
1-Phenoxypropan...	11.587	3.5	ng/ul	50705	2	10.853	287526	20.0
Phenol, p-tert-...	12.193	3.1	ng/ul	45188	2	10.853	287526	20.0
Naphthalene, 2,...	13.996	2.9	ng/ul	60572	3	14.678	414440	20.0
unknown-02	15.130	3.3	ng/ul	68710	3	14.678	414440	20.0
Diethyltoluamide	15.406	3.8	ng/ul	78008	3	14.678	414440	20.0