

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
 Data File : BG063071.D
 Acq On : 27 Sep 2024 13:13
 Operator : RC/JU
 Sample : P3927-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDCC7

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-BG092524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063071.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.317	37	39	43	rVB5	10520	11600	0.48%	0.044%
2	3.722	104	108	114	rBV	69433	113850	4.74%	0.429%
3	5.979	488	492	496	rVB4	5006	8164	0.34%	0.031%
4	6.684	610	612	616	rVV2	5679	9029	0.38%	0.034%
5	6.737	617	621	622	rVV2	6655	8280	0.34%	0.031%
6	6.772	625	627	630	rVV2	5013	6352	0.26%	0.024%
7	6.807	630	633	636	rVV3	7192	8824	0.37%	0.033%
8	6.848	636	640	646	rVV6	12205	22476	0.94%	0.085%
9	6.907	646	650	652	rVV3	8684	10839	0.45%	0.041%
10	6.936	652	655	656	rVV2	6060	6803	0.28%	0.026%
11	7.019	663	669	678	rVV2	153275	278506	11.59%	1.051%
12	7.083	678	680	686	rVB5	6815	9082	0.38%	0.034%
13	7.177	691	696	702	rBV	93876	155704	6.48%	0.587%
14	7.312	717	719	721	rBV2	6531	6617	0.28%	0.025%
15	7.383	725	731	740	rVB2	159595	288343	12.00%	1.088%
16	7.495	747	750	751	rBV3	6957	6398	0.27%	0.024%
17	7.665	774	779	780	rBV4	6557	9920	0.41%	0.037%
18	7.747	788	793	796	rBV5	8341	14212	0.59%	0.054%
19	7.771	796	797	799	rVV2	9870	7415	0.31%	0.028%
20	7.847	801	810	818	rVV	247417	436832	18.18%	1.648%
21	7.912	818	821	829	rVV3	21076	40957	1.70%	0.154%
22	8.041	840	843	847	rBV5	11014	14430	0.60%	0.054%
23	8.276	881	883	886	rBV3	5766	6938	0.29%	0.026%
24	8.382	898	901	905	rVV5	9107	14844	0.62%	0.056%
25	8.440	909	911	915	rVB3	9122	9902	0.41%	0.037%
26	8.552	925	930	938	rVB	217928	362018	15.07%	1.366%
27	8.611	938	940	943	rBV4	11407	12735	0.53%	0.048%
28	8.799	968	972	974	rBV3	5514	9650	0.40%	0.036%
29	8.834	974	978	985	rVB4	45859	72065	3.00%	0.272%
30	8.963	990	1000	1002	rBV6	32325	69424	2.89%	0.262%
31	8.999	1002	1006	1013	rVV2	137261	229165	9.54%	0.864%
32	9.081	1017	1020	1024	rVB3	21321	25418	1.06%	0.096%
33	9.163	1032	1034	1036	rBV2	7779	6571	0.27%	0.025%
34	9.404	1072	1075	1077	rBV4	5627	6809	0.28%	0.026%
35	9.428	1077	1079	1084	rVB4	10159	13380	0.56%	0.050%
36	9.504	1086	1092	1093	rBV4	7676	13149	0.55%	0.050%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
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37	9.563	1101	1102	1107	rVB3	13942	12003	0.50%	0.045%
38	9.633	1107	1114	1120	rVB6	14897	35113	1.46%	0.132%
39	9.721	1120	1129	1135	rBV2	140950	272165	11.33%	1.027%
40	9.927	1162	1164	1167	rVB2	8429	8338	0.35%	0.031%

41	9.992	1173	1175	1181	rVB5	8386	9889	0.41%	0.037%
42	10.050	1181	1185	1186	rBV3	6000	6803	0.28%	0.026%
43	10.180	1201	1207	1211	rVB7	8772	16240	0.68%	0.061%
44	10.221	1211	1214	1215	rBV3	8244	7046	0.29%	0.027%
45	10.262	1215	1221	1231	rVB	254569	452978	18.85%	1.709%

46	10.632	1277	1284	1291	rVV2	888477	1528401	63.60%	5.765%
47	10.755	1297	1305	1320	rBV2	303464	785700	32.70%	2.964%
48	10.932	1329	1335	1339	rVB5	23286	38473	1.60%	0.145%
49	11.032	1347	1352	1358	rBV8	23723	40927	1.70%	0.154%
50	11.114	1360	1366	1369	rBV3	22118	37033	1.54%	0.140%

51	11.255	1385	1390	1395	rVB7	17959	28443	1.18%	0.107%
52	11.378	1409	1411	1414	rBV4	9330	8849	0.37%	0.033%
53	11.455	1420	1424	1425	rBV4	17061	15990	0.67%	0.060%
54	11.531	1431	1437	1440	rBV2	105988	198399	8.26%	0.748%
55	11.619	1450	1452	1456	rBV5	11997	13592	0.57%	0.051%

56	11.672	1456	1461	1465	rBV3	29188	46977	1.95%	0.177%
57	11.790	1479	1481	1484	rVB5	9344	7760	0.32%	0.029%
58	11.878	1491	1496	1501	rBV3	56485	86461	3.60%	0.326%
59	11.925	1502	1504	1506	rVB3	9136	6316	0.26%	0.024%
60	12.054	1520	1526	1532	rBV4	89207	169734	7.06%	0.640%

61	12.142	1538	1541	1546	rVB6	14682	21267	0.89%	0.080%
62	12.424	1587	1589	1593	rVB5	7218	7454	0.31%	0.028%
63	12.477	1596	1598	1599	rBV2	10781	10160	0.42%	0.038%
64	12.489	1599	1600	1602	rVB2	14273	7833	0.33%	0.030%
65	12.806	1652	1654	1658	rBV5	7238	7710	0.32%	0.029%

66	13.023	1688	1691	1693	rBV4	22374	24846	1.03%	0.094%
67	13.141	1709	1711	1713	rBV3	12253	9261	0.39%	0.035%
68	13.311	1736	1740	1746	rVB2	49308	76168	3.17%	0.287%
69	13.399	1754	1755	1757	rBV2	11423	10478	0.44%	0.040%
70	13.652	1794	1798	1803	rVB	100046	130336	5.42%	0.492%

71	13.770	1816	1818	1821	rBV4	18289	20709	0.86%	0.078%
72	13.887	1833	1838	1845	rBV	589347	878479	36.56%	3.314%
73	13.975	1849	1853	1856	rVB4	41901	58091	2.42%	0.219%
74	14.028	1856	1862	1866	rBV6	54497	101988	4.24%	0.385%
75	14.175	1881	1887	1895	rVB	645727	995290	41.42%	3.754%

76	14.257	1898	1901	1906	rVV6	47462	58235	2.42%	0.220%
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77	14.386	1918	1923	1927	rBV4	94857	152693	6.35%	0.576%
78	14.428	1927	1930	1934	rVV4	87128	138269	5.75%	0.522%
79	14.486	1934	1940	1947	rVB	680020	1159131	48.24%	4.372%
80	14.692	1970	1975	1979	rBV	191616	314883	13.10%	1.188%
81	14.733	1979	1982	1986	rVB3	102290	131708	5.48%	0.497%
82	14.827	1995	1998	2005	rVB6	61149	99340	4.13%	0.375%
83	15.338	2083	2085	2091	rVB4	73903	79175	3.29%	0.299%
84	15.479	2103	2109	2115	rVB	953087	1414988	58.88%	5.337%
85	15.603	2124	2130	2143	rVB2	236451	648207	26.97%	2.445%
86	15.844	2166	2171	2176	rVB	289632	426319	17.74%	1.608%
87	16.208	2229	2233	2244	rVB4	115288	256576	10.68%	0.968%
88	16.889	2345	2349	2353	rBV2	121560	172824	7.19%	0.652%
89	16.995	2365	2367	2371	rVB2	74694	74594	3.10%	0.281%
90	17.236	2402	2408	2414	rBV	1060046	1597095	66.46%	6.024%
91	17.330	2419	2424	2430	rVB2	1186755	1748707	72.77%	6.596%
92	17.735	2490	2493	2497	rVB2	80206	98670	4.11%	0.372%
93	18.029	2541	2543	2549	rVB6	68544	81079	3.37%	0.306%
94	18.129	2557	2560	2564	rBV3	100199	134498	5.60%	0.507%
95	18.429	2609	2611	2616	rVB5	70955	72248	3.01%	0.273%
96	19.627	2810	2815	2821	rBV	1598158	2325961	96.79%	8.773%
97	21.484	3126	3131	3138	rVB2	1269878	1957170	81.45%	7.382%
98	22.248	3258	3261	3266	rVB2	125750	187983	7.82%	0.709%
99	24.316	3604	3613	3625	rVB	946975	2403025	100.00%	9.064%
100	24.528	3639	3649	3661	rVB	836133	2257460	93.94%	8.515%

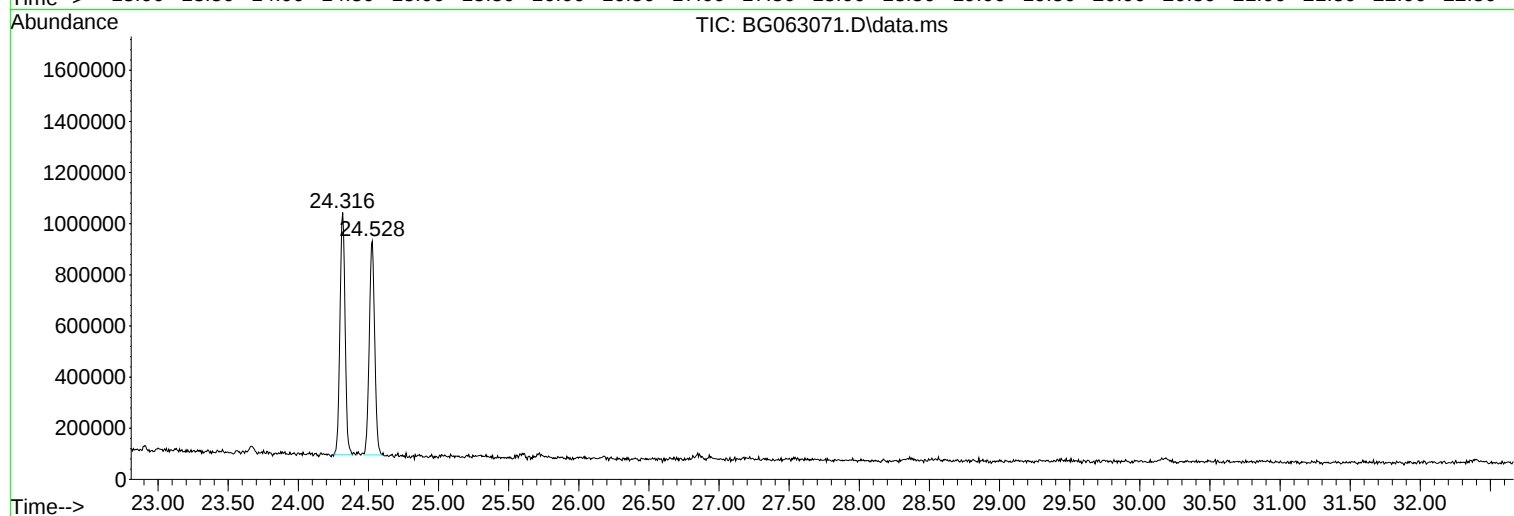
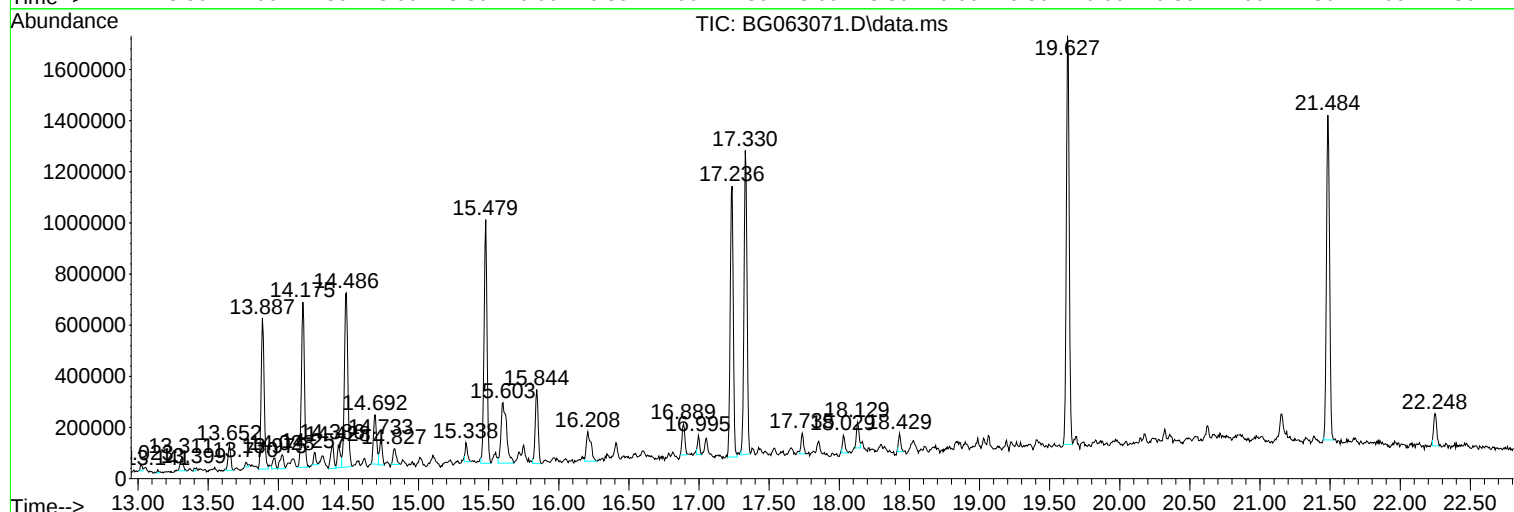
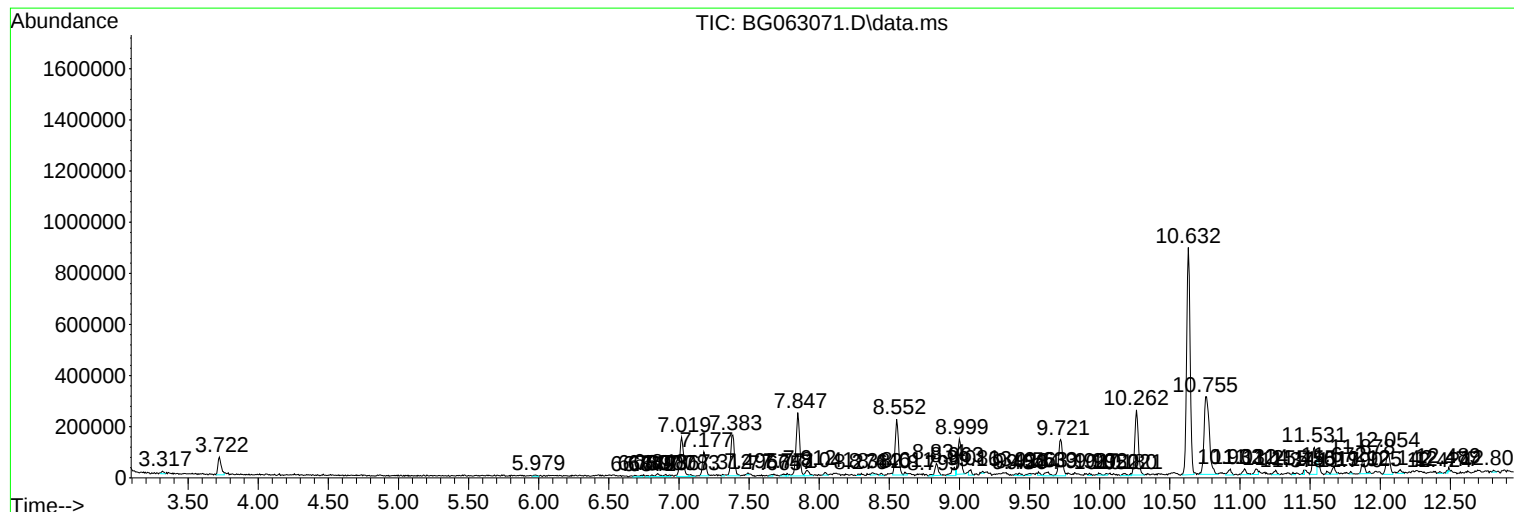
Sum of corrected areas: 26511239

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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCC7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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



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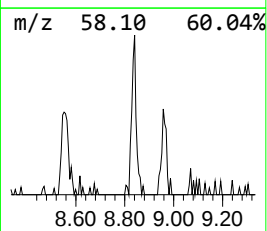
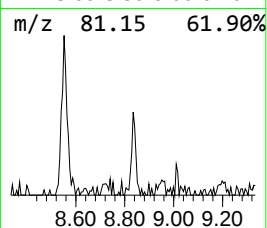
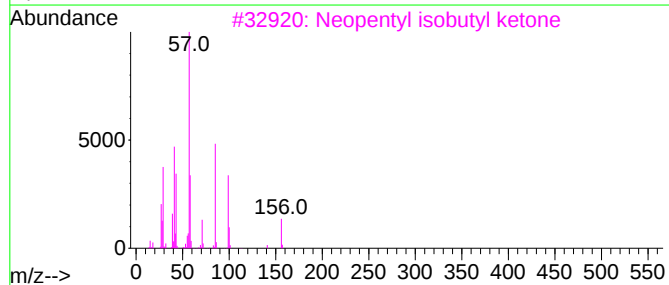
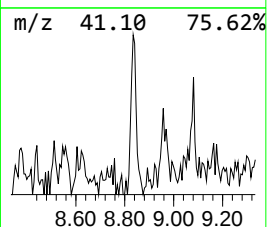
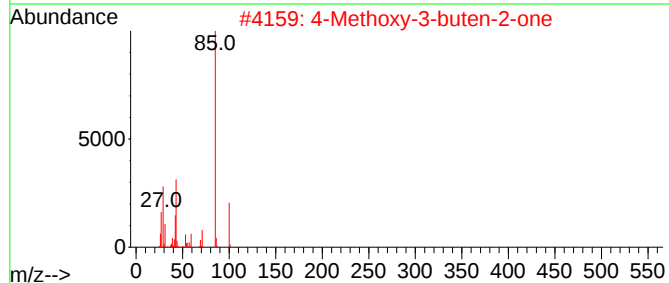
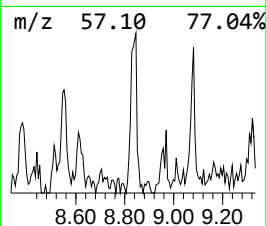
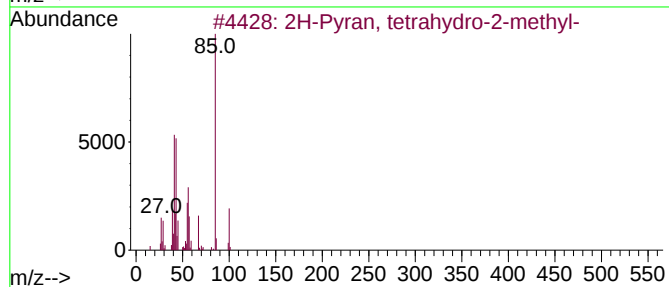
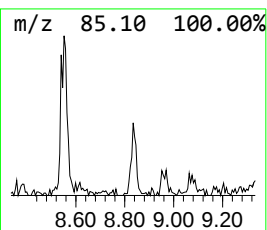
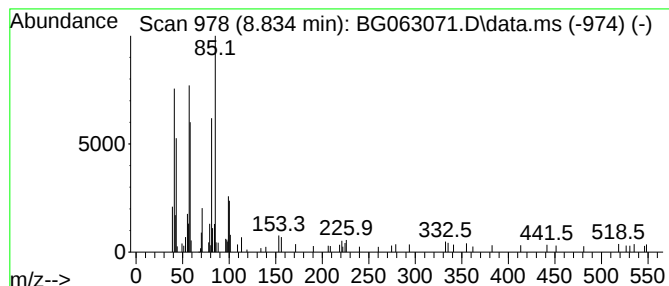
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.834	3.30 ng/ul	72065	1,4-Dichlorobenzene-d4	7.847

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	53
2	4-Methoxy-3-buten-2-one			100	C5H8O2	004652-27-1	50
3	Neopentyl isobutyl ketone			156	C10H20O	040239-19-8	50
4	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	47
5	Pentanoic acid, 2-propenyl ester			142	C8H14O2	006321-45-5	47



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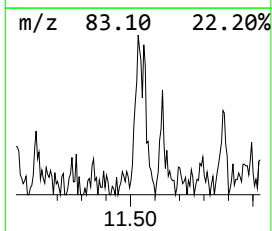
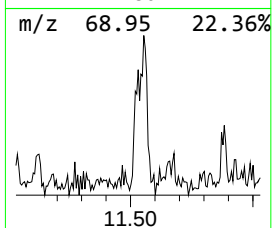
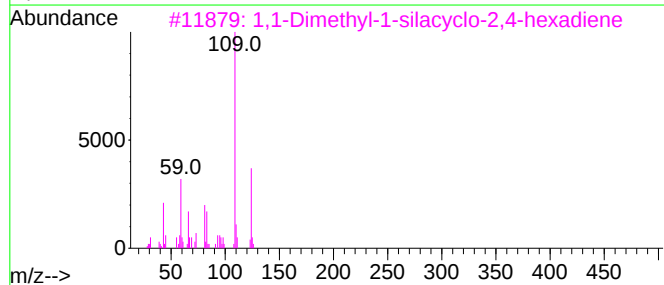
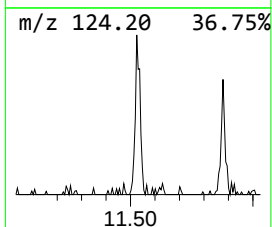
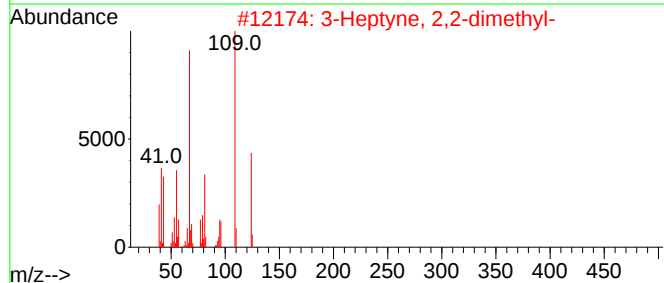
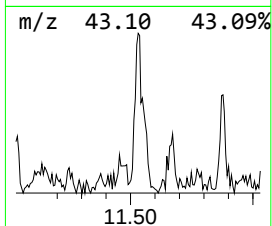
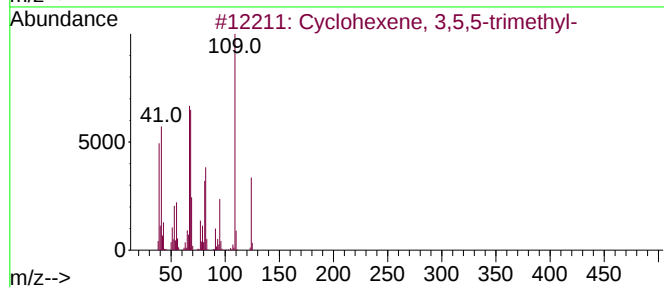
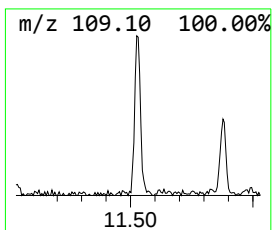
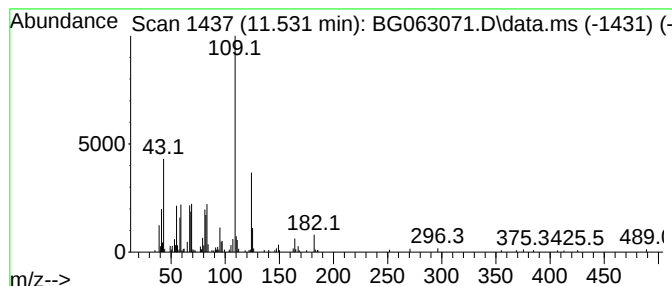
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclohexene, 3,5,5-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.531	2.60 ng/ul	198399	Naphthalene-d8	10.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3,5,5-trimethyl-	124	C9H16	000933-12-0	62
2			3-Heptyne, 2,2-dimethyl-	124	C9H16	029022-29-5	53
3			1,1-Dimethyl-1-silacyclo-2,4-hex...	124	C7H12Si	052023-18-4	53
4			1,3-Heptadiene, 2,3-dimethyl-	124	C9H16	074779-65-0	52
5			p-Fluoroethylbenzene	124	C8H9F	000459-47-2	50



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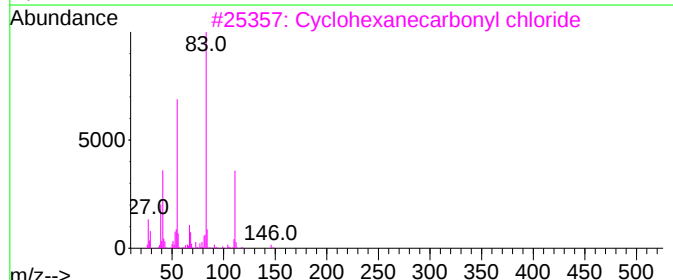
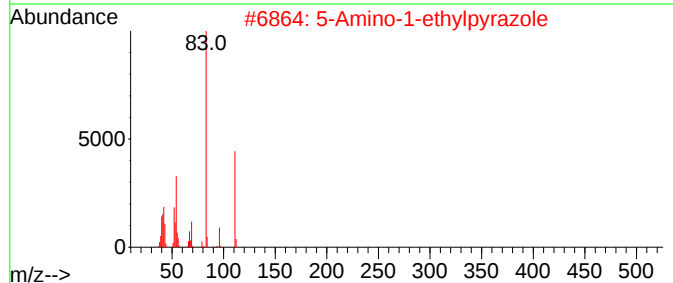
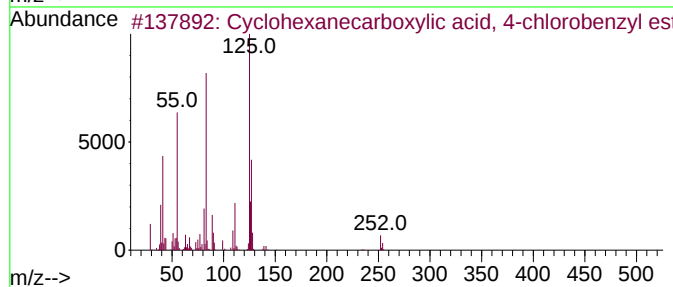
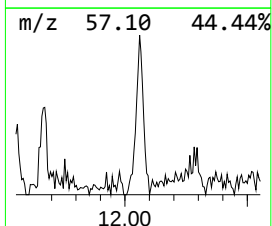
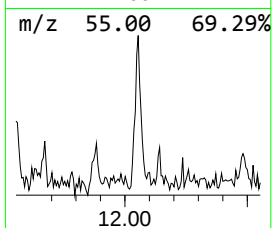
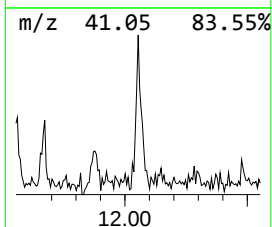
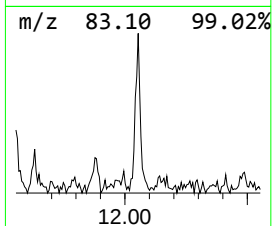
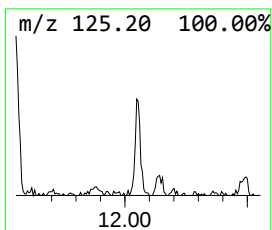
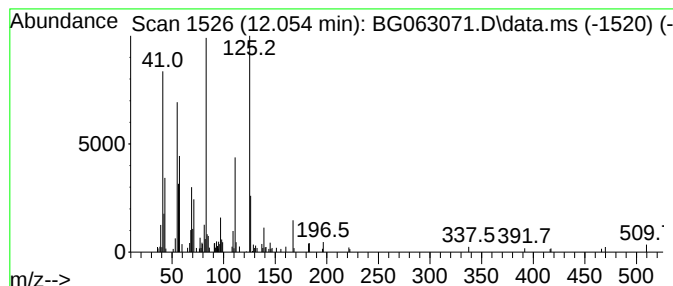
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexanecarboxylic acid,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.054	2.22 ng/ul	169734	Naphthalene-d8	10.638

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanecarboxylic acid, 4-ch...	252	C14H17ClO2	1000279-91-6	72
2			5-Amino-1-ethylpyrazole	111	C5H9N3	003528-58-3	43
3			Cyclohexanecarbonyl chloride	146	C7H11ClO	002719-27-9	43
4			Cyclohexane carboxylic acid hydr...	142	C7H14N2O	038941-47-8	38
5			3-Cyclopentylpropionic acid, 3-p...	352	C23H44O2	1000293-57-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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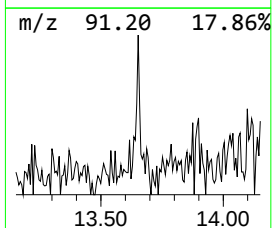
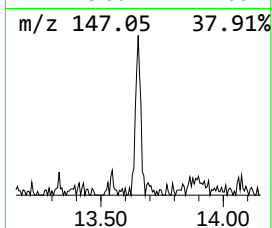
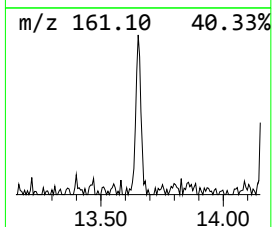
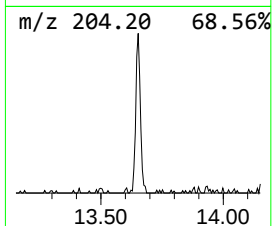
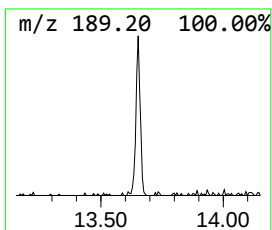
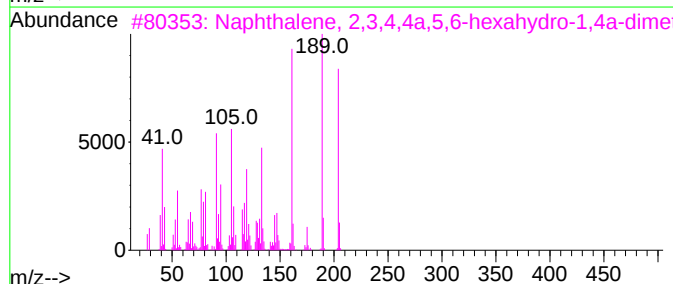
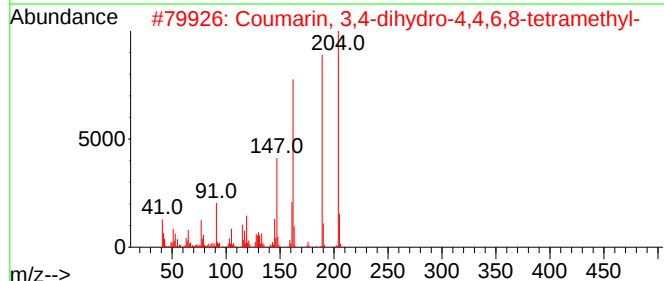
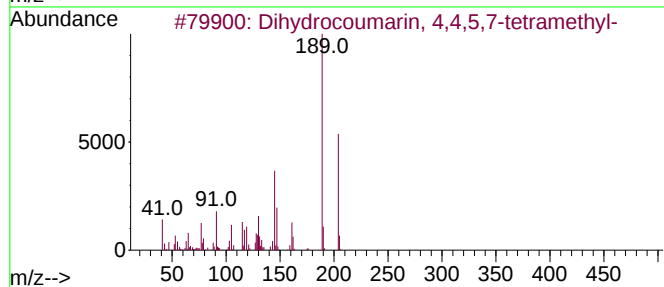
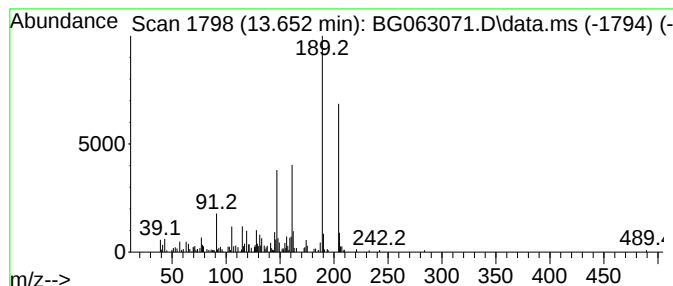
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dihydrocoumarin, 4,4,5,7-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.652	2.25 ng/ul	130336	Acenaphthene-d10	14.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dihydrocoumarin, 4,4,5,7-tetrame...	204	C13H16O2	040662-14-4	64
2			Coumarin, 3,4-dihydro-4,4,6,8-te...	204	C13H16O2	249629-60-5	60
3			Naphthalene, 2,3,4,4a,5,6-hexahy...	204	C15H24	000473-14-3	58
4			3,4-Dihydrocoumarin, 4,4-dimethy...	204	C13H16O2	029423-74-3	58
5			Benzene, 1,3,5-tris(1-methylethyl)-	204	C15H24	000717-74-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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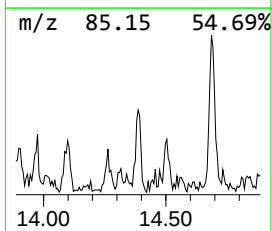
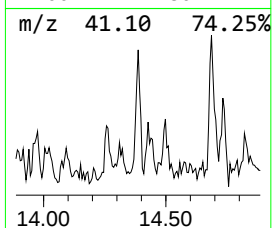
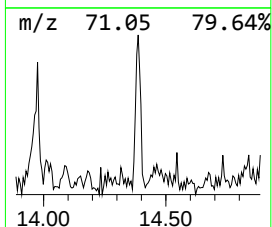
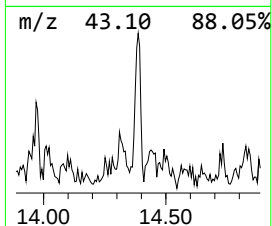
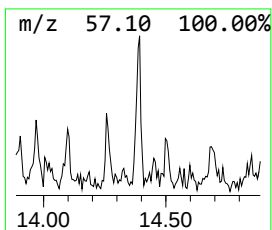
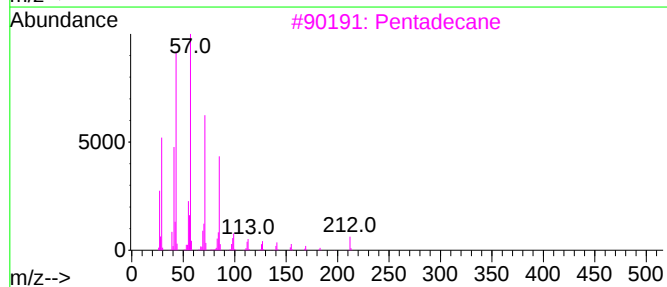
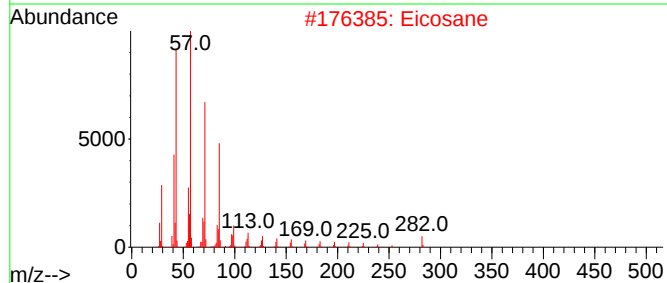
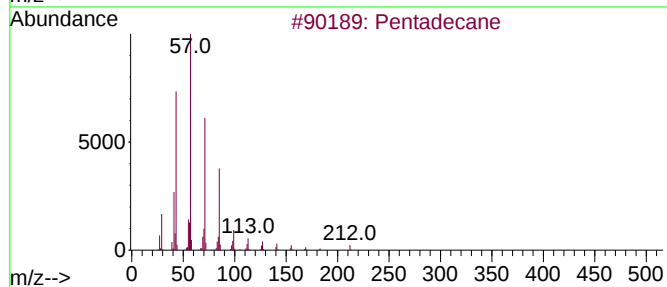
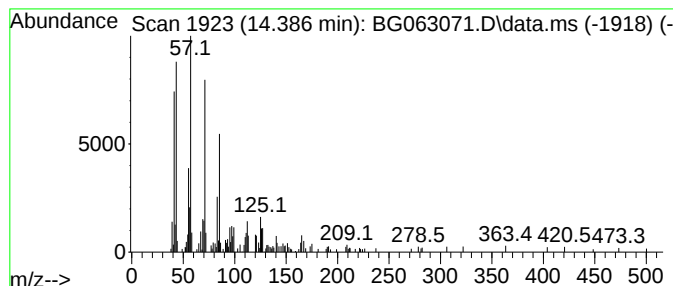
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.387	2.63 ng/ul	152693	Acenaphthene-d10	14.480

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	90
2		Eicosane	282	C20H42	000112-95-8	90
3		Pentadecane	212	C15H32	000629-62-9	87
4		Pentadecane	212	C15H32	000629-62-9	83
5		Hexadecane	226	C16H34	000544-76-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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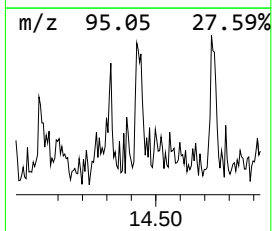
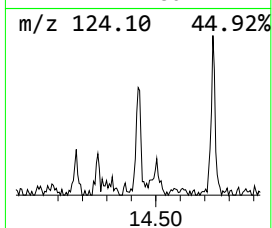
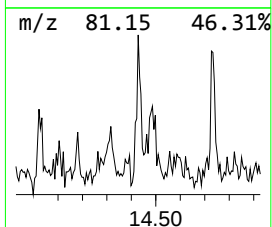
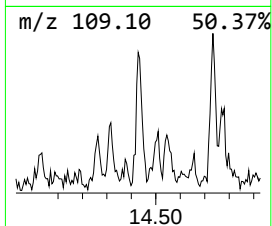
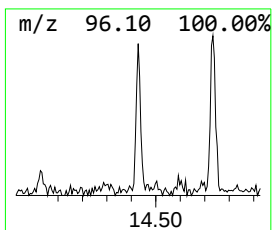
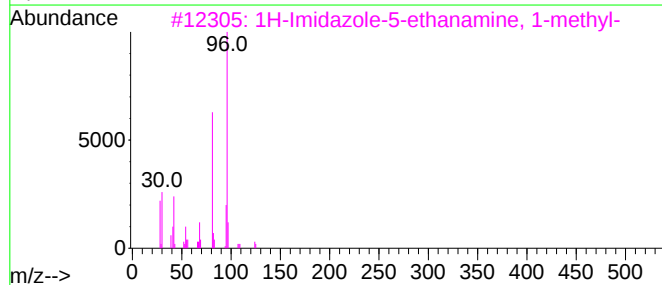
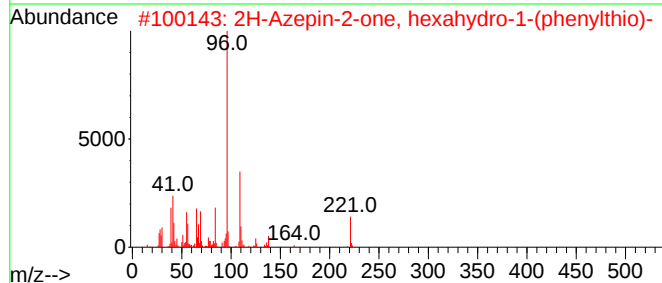
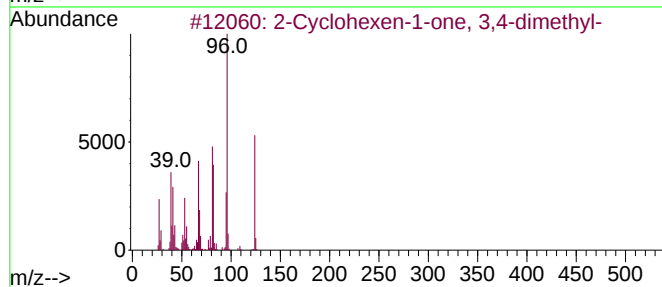
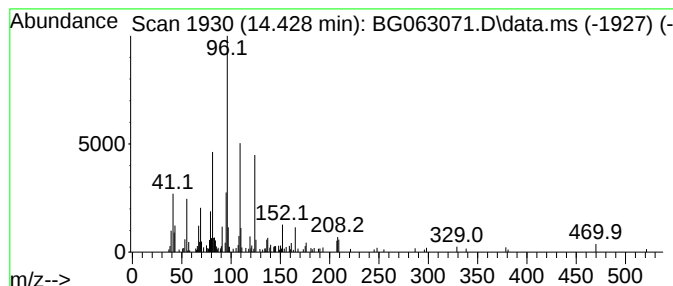
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Cyclohexen-1-one, 3,4-dim... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.428	2.39 ng/ul	138269	Acenaphthene-d10	14.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one, 3,4-dimethyl-	124	C8H12O	1000197-00-4	52
2			2H-Azepin-2-one, hexahydro-1-(ph...	221	C12H15NOS	036452-23-0	43
3			1H-Imidazole-5-ethanamine, 1-met...	125	C6H11N3	000644-42-8	38
4			Ethylamine, 1-methyl-2-(5-methyl...	139	C7H13N3	1000316-41-3	35
5			1H-Imidazole-4-ethanamine, 2-met...	125	C6H11N3	034392-54-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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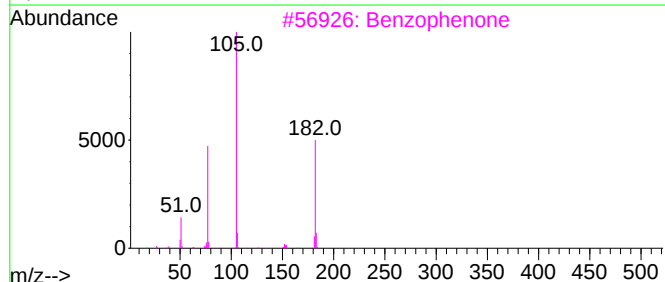
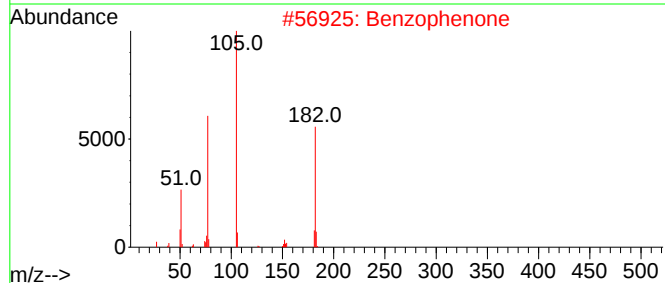
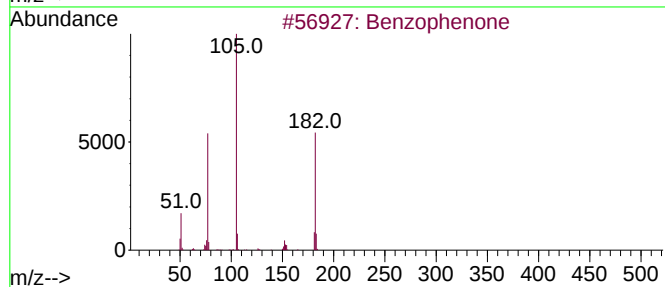
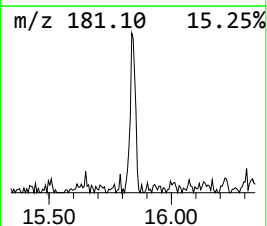
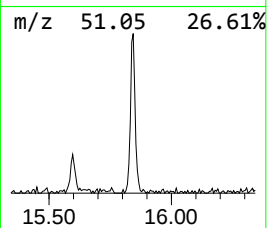
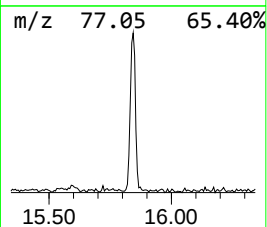
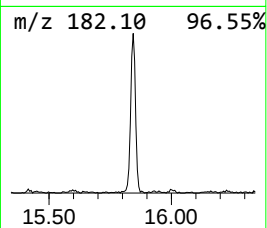
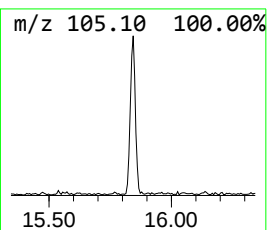
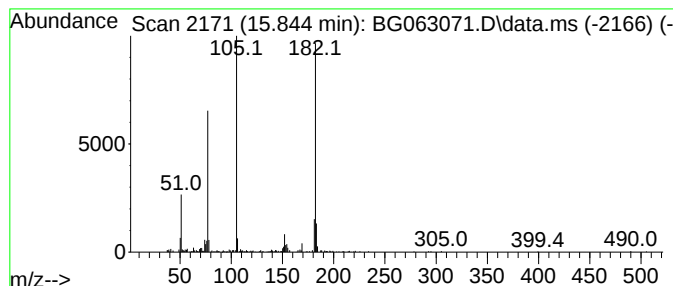
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.844	7.36 ng/ul	426319	Acenaphthene-d10	14.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	93
2			Benzophenone	182	C13H10O	000119-61-9	93
3			Benzophenone	182	C13H10O	000119-61-9	90
4			Benzophenone	182	C13H10O	000119-61-9	86
5			Benzoic acid, 3,4-dimethoxy-	182	C9H10O4	000093-07-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

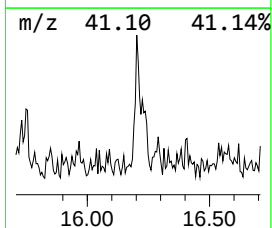
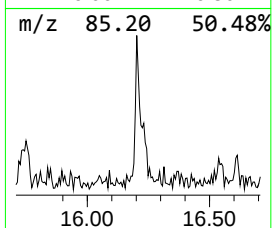
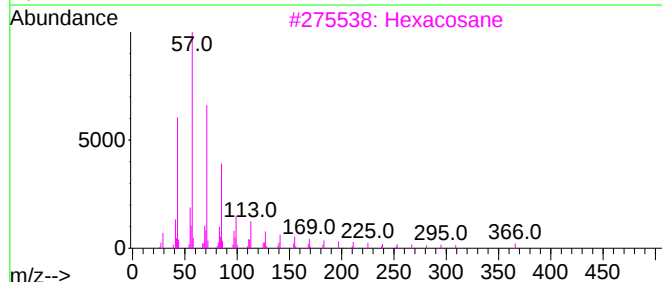
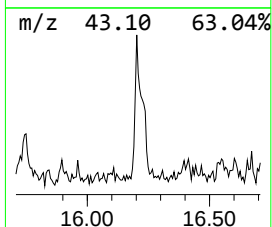
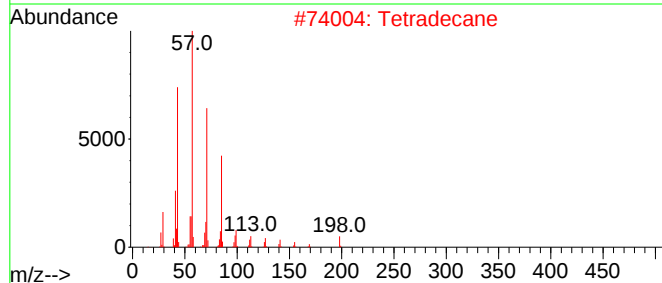
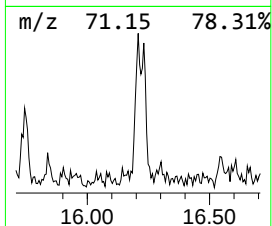
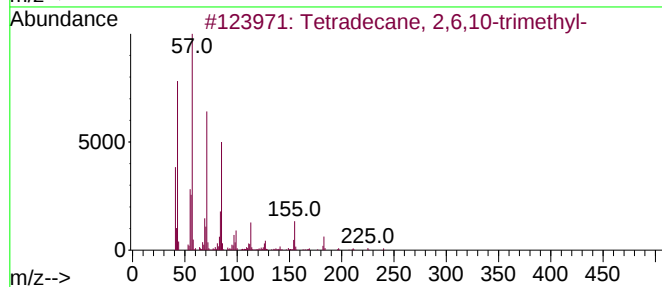
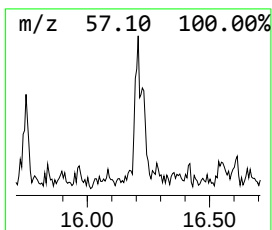
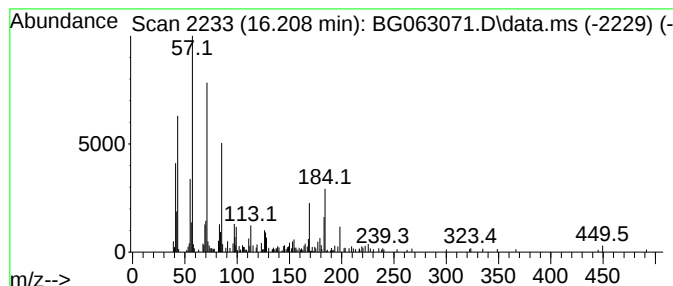
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.208	3.21 ng/ul	256576	Phenanthrene-d10		17.236	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	83
2	Tetradecane		198	C14H30	000629-59-4	70
3	Hexacosane		366	C26H54	000630-01-3	64
4	Tetradecane		198	C14H30	000629-59-4	64
5	Tetradecane		198	C14H30	000629-59-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063071.D
Acq On : 27 Sep 2024 13:13
Operator : RC/JU
Sample : P3927-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCC7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.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
2H-Pyran, tetra...	8.834	3.3	ng/ul	72065	1	7.847	436832	20.0
Cyclohexene, 3,...	11.531	2.6	ng/ul	198399	2	10.638	1528400	20.0
Cyclohexanecarb...	12.054	2.2	ng/ul	169734	2	10.638	1528400	20.0
Dihydrocoumarin...	13.652	2.3	ng/ul	130336	3	14.480	1159130	20.0
(DEL) Alkane: S...	14.387	2.6	ng/ul	152693	3	14.480	1159130	20.0
2-Cyclohexen-1-...	14.428	2.4	ng/ul	138269	3	14.480	1159130	20.0
Benzophenone	15.844	7.4	ng/ul	426319	3	14.480	1159130	20.0
(DEL) Alkane: S...	16.208	3.2	ng/ul	256576	4	17.236	1597100	20.0