

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
 Data File : BG061847.D
 Acq On : 2 Jul 2024 15:04
 Operator : MA/JU
 Sample : P2963-15
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CR9

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-BG062024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061847.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.467	25	30	32	rBV3	12233	17176	0.96%	0.072%
2	3.731	71	75	86	rVB	25070	50530	2.83%	0.213%
3	3.895	101	103	107	rBV4	6455	9241	0.52%	0.039%
4	3.989	115	119	125	rVV3	18387	29194	1.63%	0.123%
5	4.060	127	131	136	rBV4	4694	7976	0.45%	0.034%
6	4.160	143	148	149	rBV4	3501	5268	0.29%	0.022%
7	4.213	153	157	161	rBV3	15080	25979	1.45%	0.109%
8	4.583	219	220	225	rVV4	7378	7861	0.44%	0.033%
9	4.730	241	245	249	rBV4	9478	18077	1.01%	0.076%
10	4.783	249	254	258	rVV3	15611	26709	1.50%	0.112%
11	4.924	276	278	283	rVB4	3699	5854	0.33%	0.025%
12	5.135	308	314	323	rVV	60662	98447	5.51%	0.415%
13	5.235	326	331	336	rBV5	6617	11887	0.67%	0.050%
14	5.376	349	355	361	rBV3	5694	11627	0.65%	0.049%
15	6.058	469	471	474	rVV4	3873	4428	0.25%	0.019%
16	6.116	476	481	489	rVB4	12295	28509	1.60%	0.120%
17	6.263	504	506	511	rVB5	3664	4788	0.27%	0.020%
18	7.039	634	638	643	rVB5	9670	15872	0.89%	0.067%
19	7.098	643	648	650	rVB5	3138	5228	0.29%	0.022%
20	7.221	662	669	681	rVB	409990	733703	41.07%	3.090%
21	7.380	690	696	705	rBV	274551	450173	25.20%	1.896%
22	7.591	725	732	739	rBV	377852	653908	36.60%	2.754%
23	7.732	754	756	761	rVB4	3766	5894	0.33%	0.025%
24	8.061	804	812	818	rBV	279702	473035	26.48%	1.992%
25	8.108	818	820	826	rVB3	12613	18014	1.01%	0.076%
26	8.437	873	876	881	rVB6	3509	5673	0.32%	0.024%
27	8.608	901	905	911	rBV8	5095	11403	0.64%	0.048%
28	8.766	924	932	940	rBV2	465710	795118	44.51%	3.349%
29	8.878	947	951	958	rVB2	27588	45911	2.57%	0.193%
30	9.225	1002	1010	1018	rBV	385362	666091	37.29%	2.806%
31	9.377	1034	1036	1039	rVV3	8819	9614	0.54%	0.040%
32	9.606	1071	1075	1084	rVB6	8414	14961	0.84%	0.063%
33	9.695	1086	1090	1092	rBV3	4685	6654	0.37%	0.028%
34	9.777	1095	1104	1109	rBV3	49536	86011	4.81%	0.362%
35	9.947	1125	1133	1144	rVV	343240	609863	34.14%	2.569%
36	10.094	1153	1158	1165	rVB7	8465	18792	1.05%	0.079%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
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37	10.159	1165	1169	1176	rBV7	9060	18052	1.01%	0.076%
38	10.300	1191	1193	1197	rVB4	4639	5015	0.28%	0.021%
39	10.488	1217	1225	1234	rBV	484540	857830	48.02%	3.613%
40	10.870	1284	1290	1297	rVV	390876	707467	39.60%	2.980%

41	10.923	1297	1299	1305	rVV	19295	24640	1.38%	0.104%
42	11.005	1307	1313	1321	rBV2	61503	114906	6.43%	0.484%
43	11.122	1331	1333	1336	rBV3	3638	5392	0.30%	0.023%
44	11.275	1358	1359	1363	rVV2	3775	4756	0.27%	0.020%
45	11.322	1365	1367	1370	rVV3	4825	5307	0.30%	0.022%

46	11.393	1373	1379	1384	rVB5	18664	33518	1.88%	0.141%
47	11.487	1393	1395	1397	rBV2	3668	4389	0.25%	0.018%
48	11.604	1409	1415	1427	rBV2	73374	146505	8.20%	0.617%
49	11.763	1437	1442	1445	rBV5	5407	10026	0.56%	0.042%
50	11.880	1458	1462	1463	rBV2	5738	5797	0.32%	0.024%

51	12.209	1516	1518	1524	rVB5	6470	7400	0.41%	0.031%
52	12.268	1524	1528	1531	rBV5	7834	10952	0.61%	0.046%
53	12.450	1554	1559	1566	rVB8	9893	17200	0.96%	0.072%
54	12.521	1568	1571	1575	rBV4	9506	14884	0.83%	0.063%
55	12.744	1601	1609	1618	rBV7	9531	26985	1.51%	0.114%

56	12.961	1640	1646	1649	rBV8	5931	11758	0.66%	0.050%
57	13.049	1659	1661	1663	rVB3	5015	4098	0.23%	0.017%
58	13.290	1699	1702	1704	rBV4	5218	6514	0.36%	0.027%
59	13.502	1733	1738	1742	rBV5	11560	22056	1.23%	0.093%
60	13.596	1752	1754	1761	rBV6	6026	11832	0.66%	0.050%

61	13.860	1797	1799	1803	rVB5	7473	7167	0.40%	0.030%
62	14.013	1821	1825	1828	rVB5	8805	13170	0.74%	0.055%
63	14.089	1831	1838	1845	rBV	793287	1133162	63.43%	4.773%
64	14.213	1854	1859	1862	rBV7	6015	11463	0.64%	0.048%
65	14.248	1862	1865	1867	rBV4	8238	8711	0.49%	0.037%

66	14.324	1874	1878	1882	rBV5	11267	17092	0.96%	0.072%
67	14.383	1882	1888	1899	rVV	830794	1354373	75.81%	5.705%
68	14.512	1908	1910	1911	rBV2	6550	4680	0.26%	0.020%
69	14.571	1918	1920	1926	rVB5	10355	11824	0.66%	0.050%
70	14.689	1933	1940	1946	rVB	591534	906042	50.72%	3.816%

71	14.753	1946	1951	1957	rVB3	28743	49959	2.80%	0.210%
72	14.888	1967	1974	1987	rBV2	337013	602761	33.74%	2.539%
73	15.000	1989	1993	1996	rBV5	7202	13360	0.75%	0.056%
74	15.088	2003	2008	2014	rVB5	25399	49434	2.77%	0.208%
75	15.141	2014	2017	2019	rBV3	8390	11502	0.64%	0.048%

76	15.341	2049	2051	2052	rBV2	5909	4136	0.23%	0.017%
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77	15.464	2069	2072	2073	rBV3	9280	10009	0.56%	0.042%
78	15.517	2078	2081	2085	rVB4	12945	17709	0.99%	0.075%
79	15.682	2102	2109	2114	rBV	1033992	1570115	87.89%	6.613%
80	15.735	2116	2118	2122	rVB3	26707	25788	1.44%	0.109%
81	15.793	2123	2128	2132	rVB2	148432	212879	11.92%	0.897%
82	15.893	2144	2145	2146	rBV	9781	5792	0.32%	0.024%
83	15.934	2150	2152	2156	rVB5	25856	28988	1.62%	0.122%
84	15.999	2158	2163	2165	rBV6	13078	24617	1.38%	0.104%
85	16.404	2225	2232	2237	rVB9	24053	59053	3.31%	0.249%
86	16.492	2245	2247	2250	rBV4	8568	10226	0.57%	0.043%
87	16.998	2331	2333	2334	rBV2	11593	7477	0.42%	0.031%
88	17.086	2344	2348	2354	rBV3	27354	40051	2.24%	0.169%
89	17.433	2401	2407	2411	rBV	671464	1027711	57.53%	4.329%
90	17.474	2411	2414	2419	rVV	401054	581783	32.57%	2.450%
91	17.532	2419	2424	2435	rVB2	1049650	1623810	90.90%	6.839%
92	18.308	2553	2556	2560	rBV2	120213	170069	9.52%	0.716%
93	18.496	2584	2588	2593	rBV3	120695	203288	11.38%	0.856%
94	18.784	2633	2637	2638	rBV4	65777	84589	4.74%	0.356%
95	18.843	2644	2647	2651	rVB	83552	116899	6.54%	0.492%
96	19.483	2751	2756	2761	rVB	871673	1237578	69.28%	5.213%
97	19.818	2807	2813	2816	rBV	1051273	1786434	100.00%	7.524%
98	21.334	3068	3071	3082	rVB4	459745	779791	43.65%	3.284%
99	21.692	3126	3132	3138	rBV2	585826	1420455	79.51%	5.983%
100	24.013	3523	3527	3533	rVB	748568	1431489	80.13%	6.029%

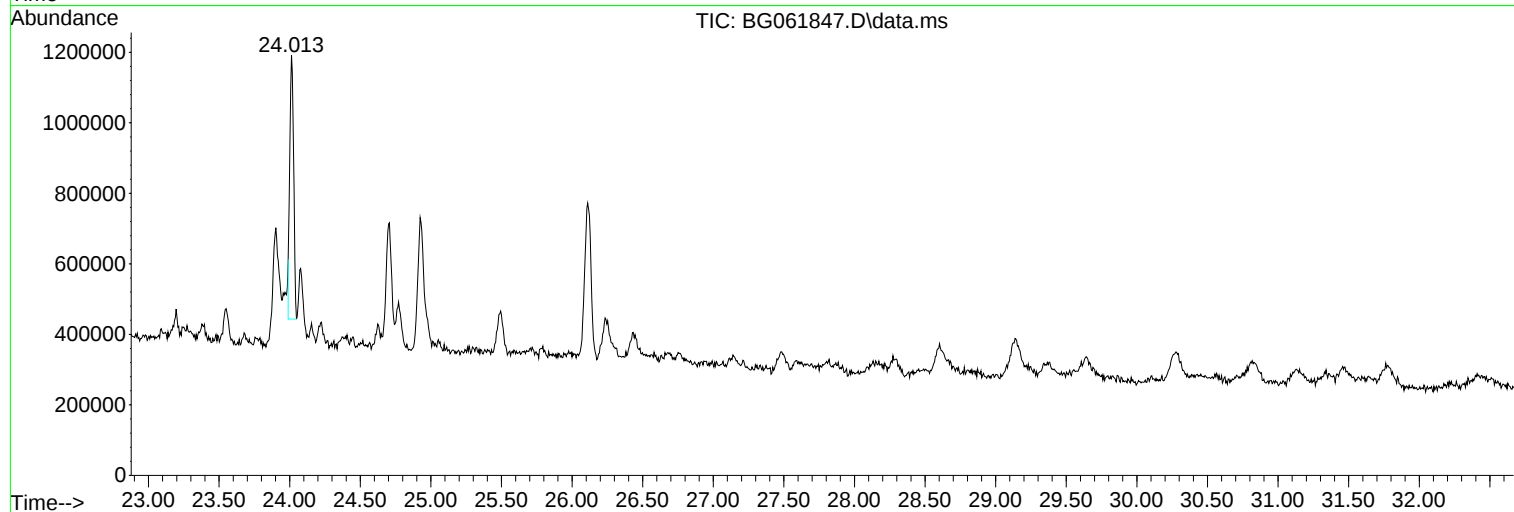
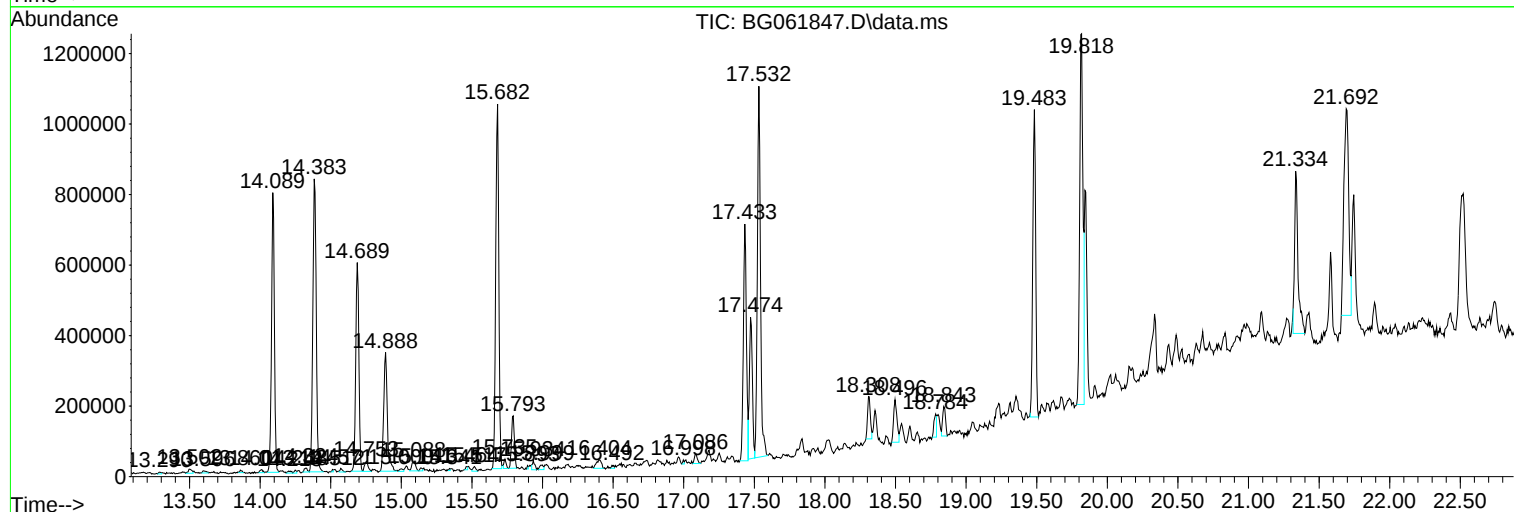
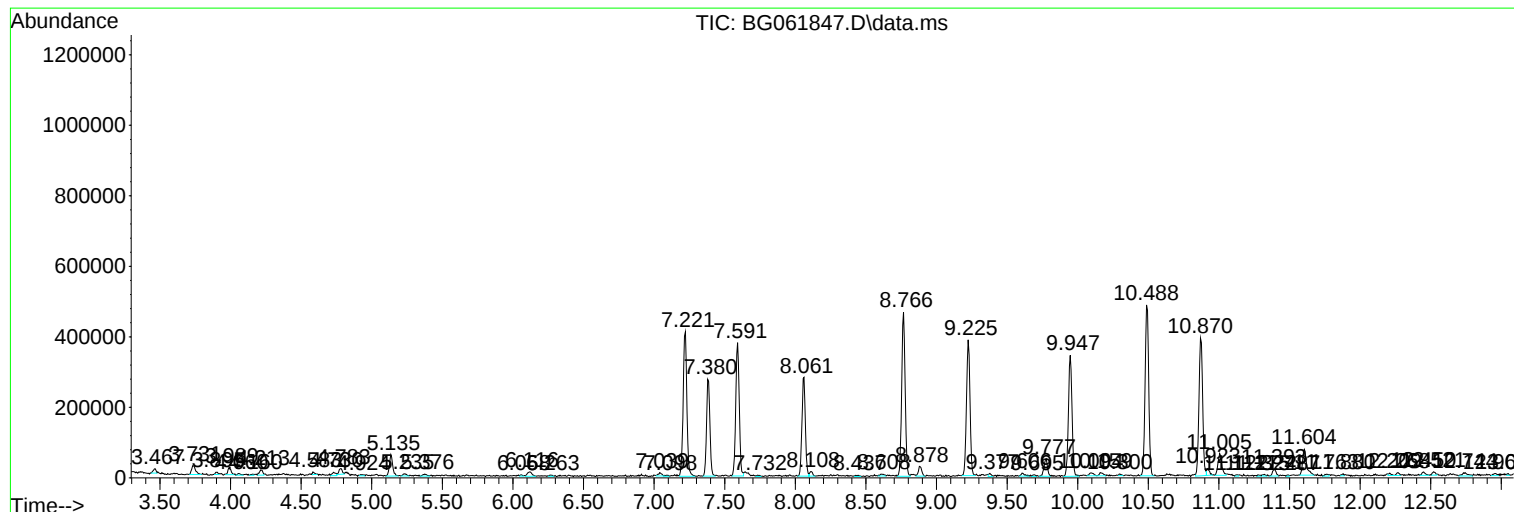
Sum of corrected areas: 23742184

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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



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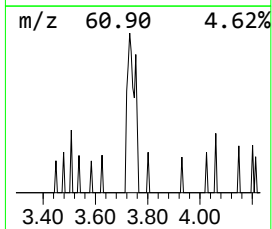
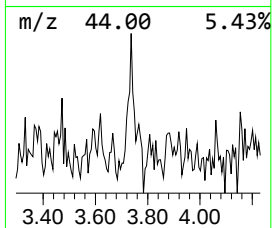
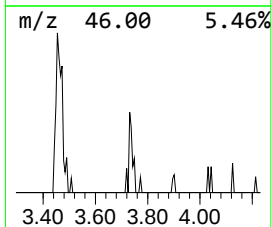
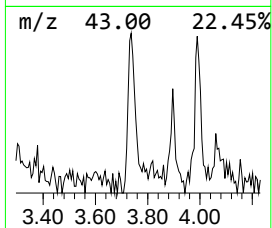
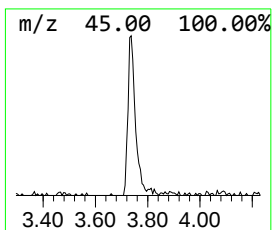
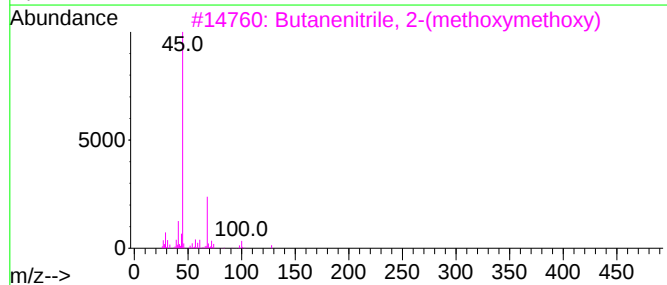
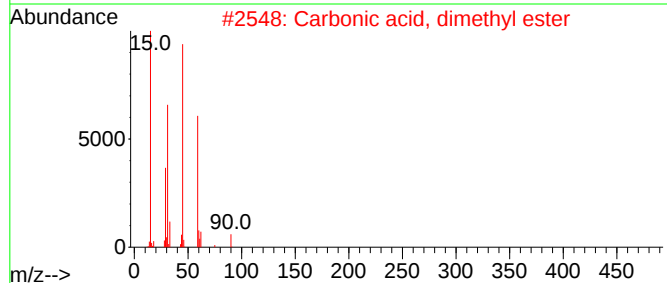
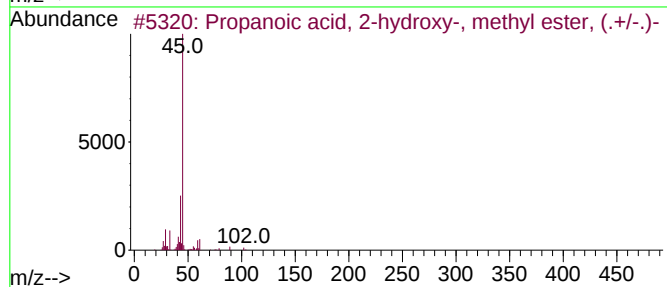
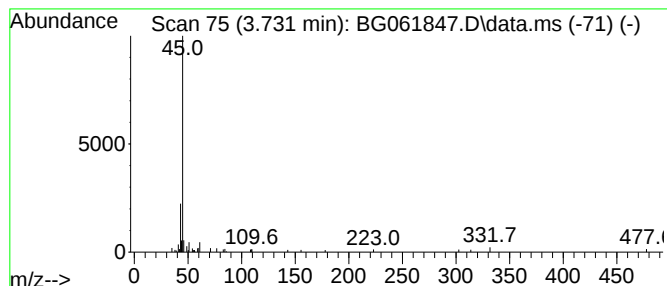
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.731	2.14 ng/ul	50530	1,4-Dichlorobenzene-d4	8.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	4
2			Carbonic acid, dimethyl ester	90	C3H6O3	000616-38-6	4
3			Butanenitrile, 2-(methoxymethoxy)	129	C6H11NO2	1000196-86-5	4
4			2-Propanol, 1,3-dimethoxy-	120	C5H12O3	000623-69-8	4
5			Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	4



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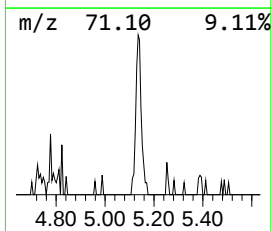
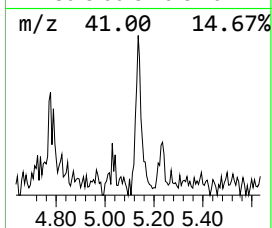
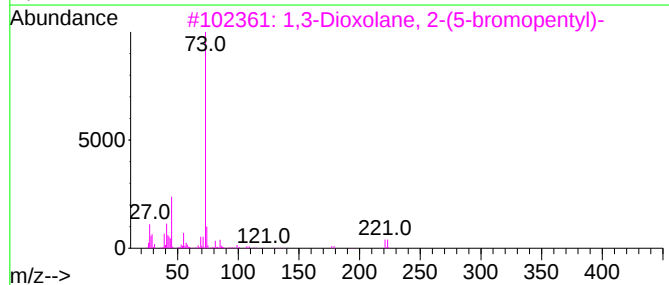
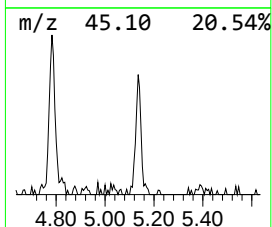
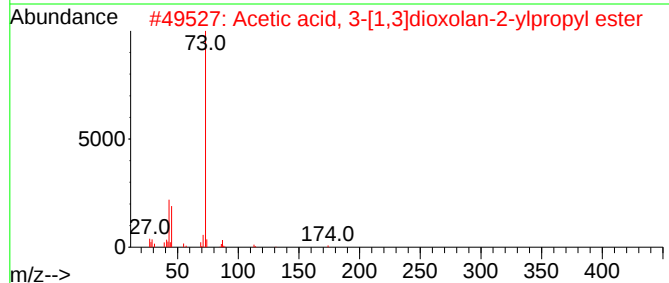
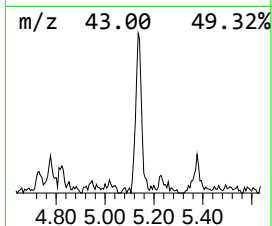
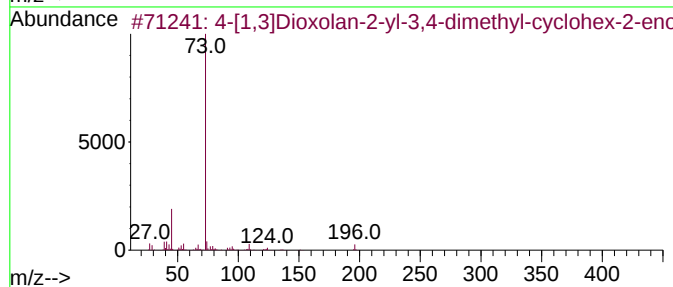
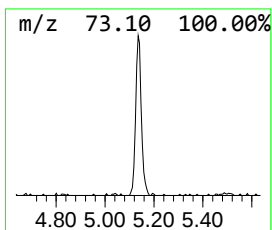
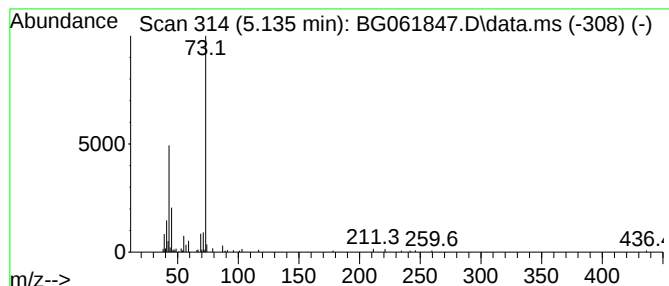
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.135	4.16 ng/ul	98447	1,4-Dichlorobenzene-d4	8.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4	-[1,3]Dioxolan-2-yl-3,4-dimethy...	196	C11H16O3	054710-16-6	37	
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	36		
3	1,3-Dioxolane, 2-(5-bromopentyl)-	222	C8H15BrO2	056741-68-5	36		
4	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	36		
5	1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	28		



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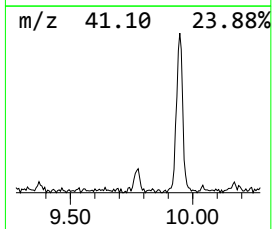
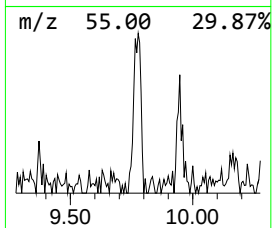
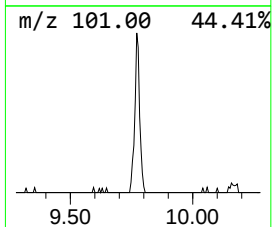
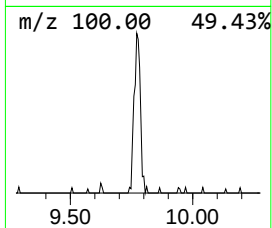
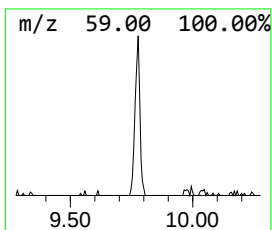
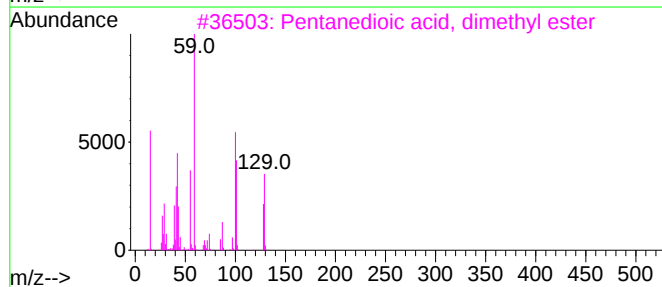
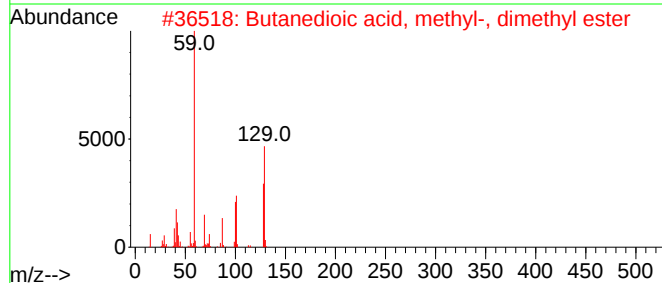
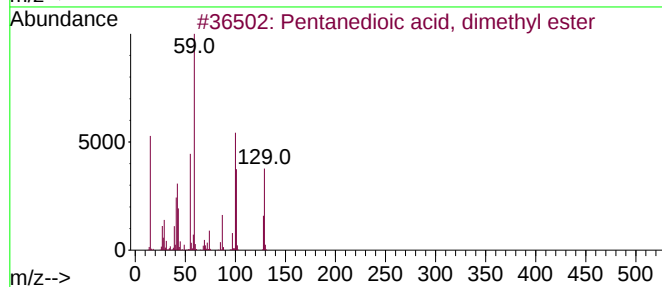
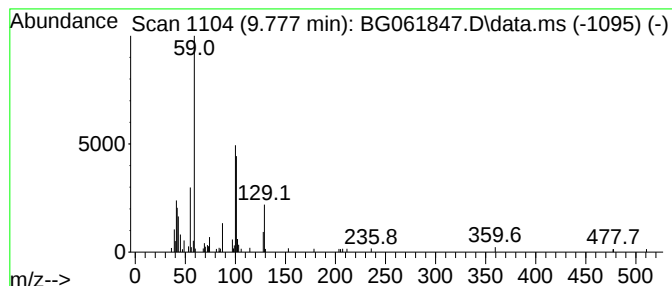
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.777	2.43 ng/ul	86011	Naphthalene-d8	10.870

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	56
2			Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	45
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	40
4			Dimethylpropylsilane	102	C5H14Si	1000432-95-4	27
5			Butanoic acid, 4-iodo-, methyl e...	228	C5H9IO2	014273-85-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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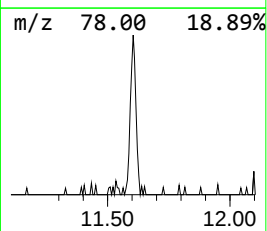
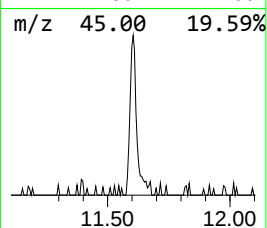
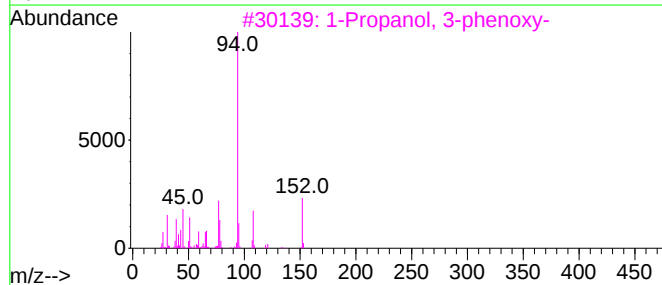
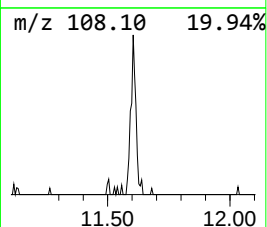
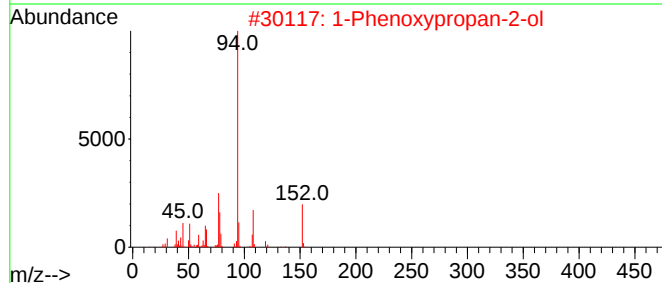
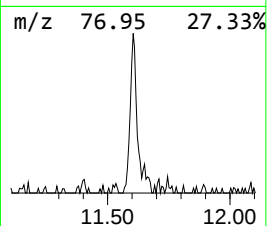
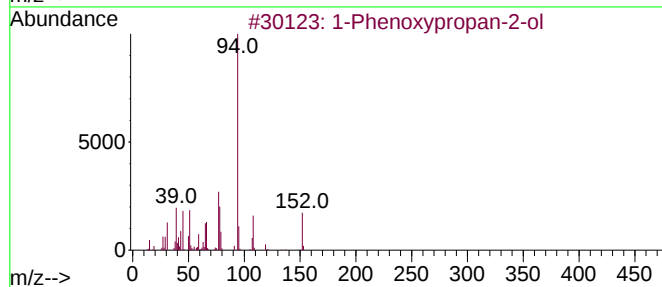
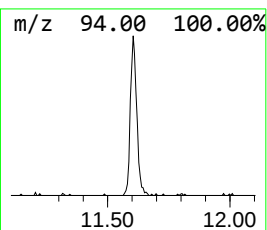
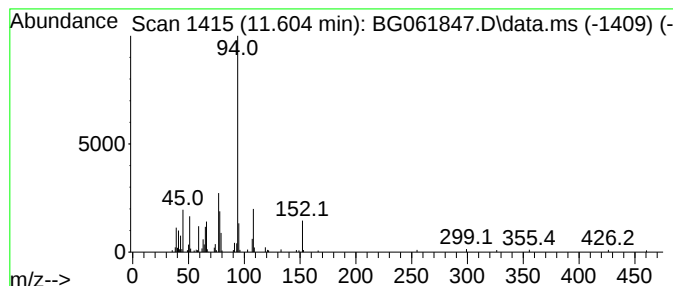
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.604	4.14 ng/ul	146505	Naphthalene-d8	10.870

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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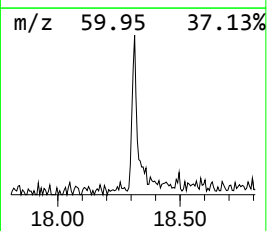
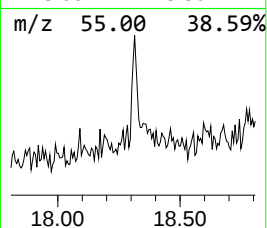
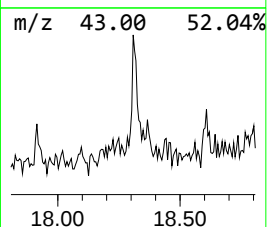
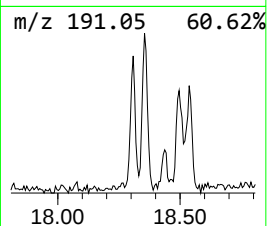
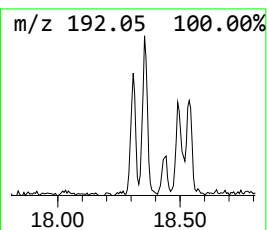
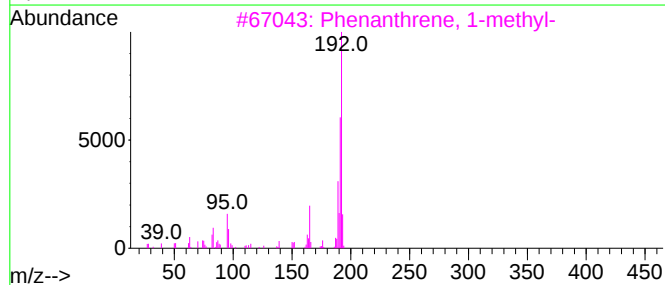
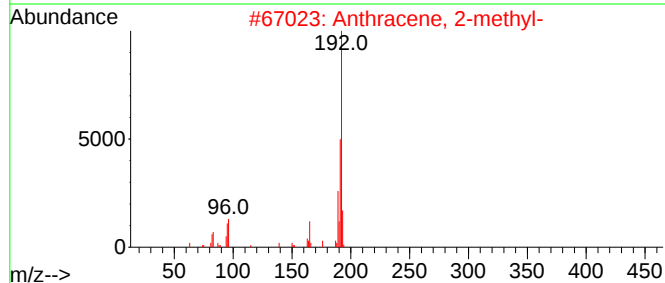
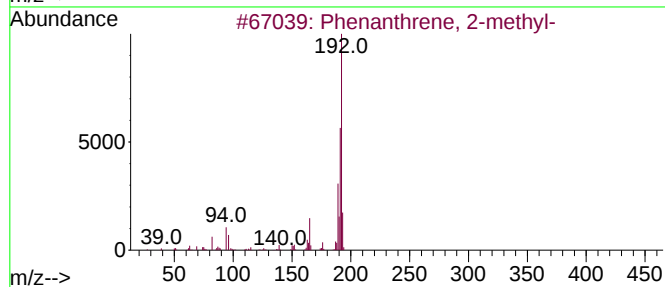
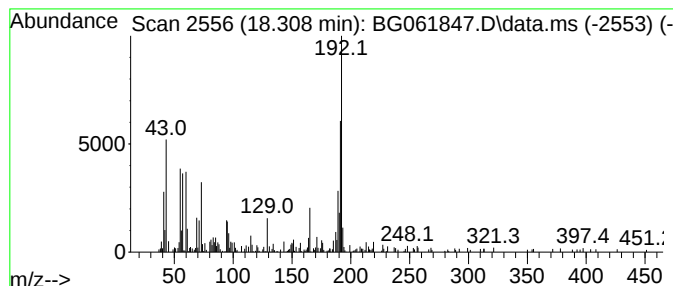
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.308	3.31 ng/ul	170069	Phenanthrene-d10	17.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	62
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	58
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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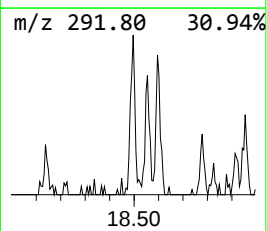
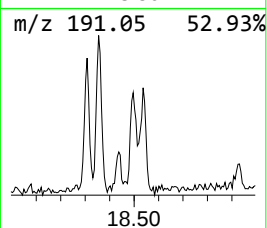
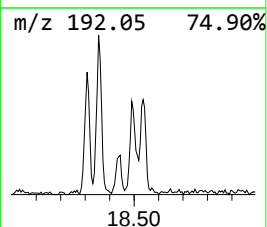
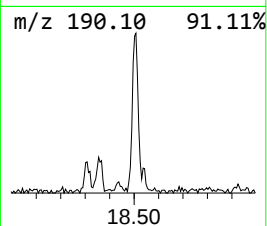
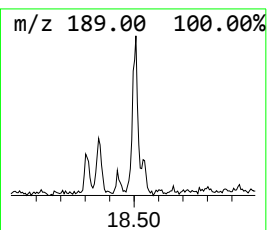
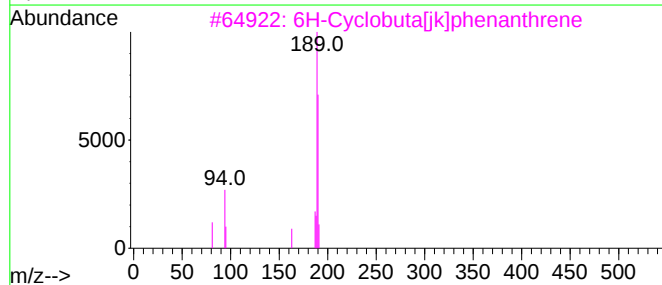
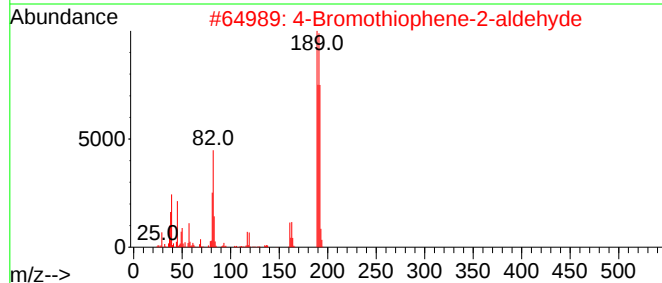
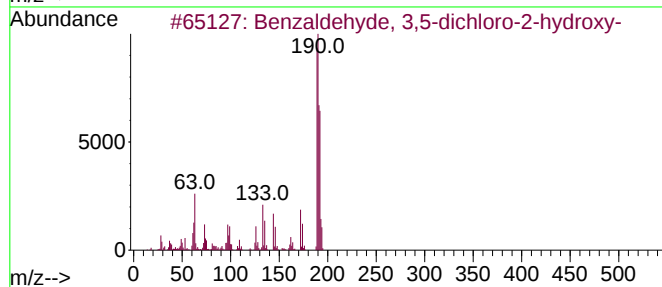
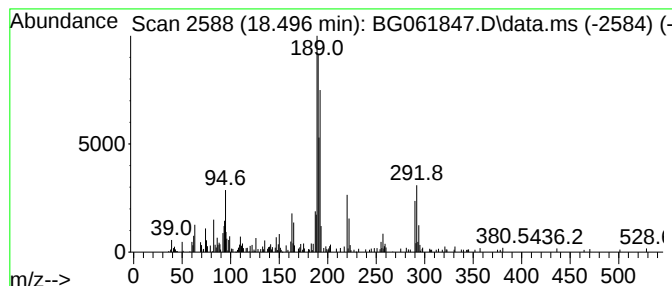
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzaldehyde, 3,5-dichloro-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.496	3.96 ng/ul	203288	Phenanthrene-d10	17.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2			4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	52
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	41
4			3-(Trifluoromethoxy)benzaldehyde	190	C8H5F3O2	052771-21-8	38
5			4-(4-Methyl-piperidin-1-yl)-phen...	190	C12H18N2	342013-25-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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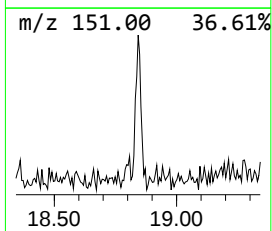
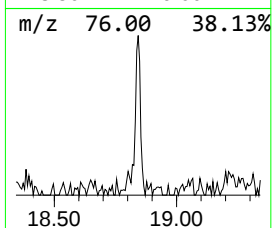
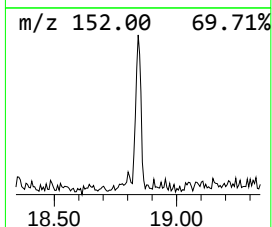
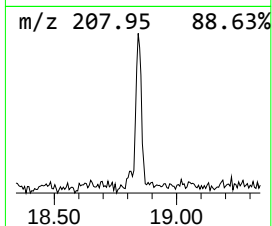
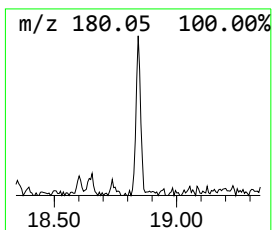
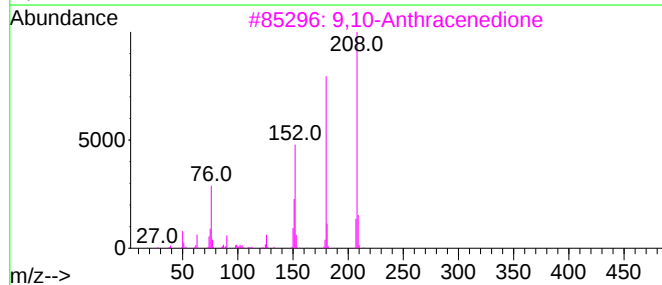
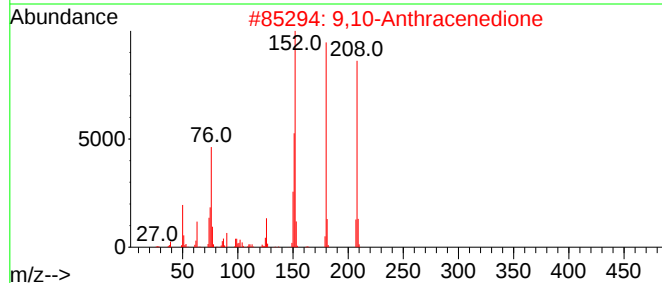
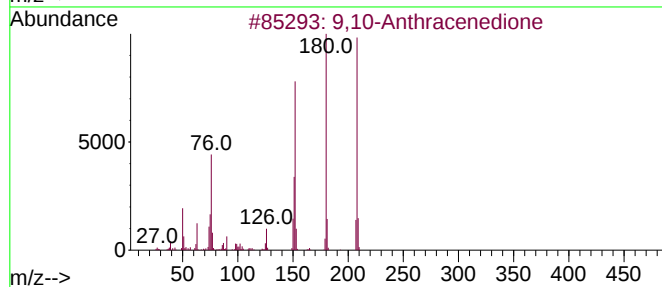
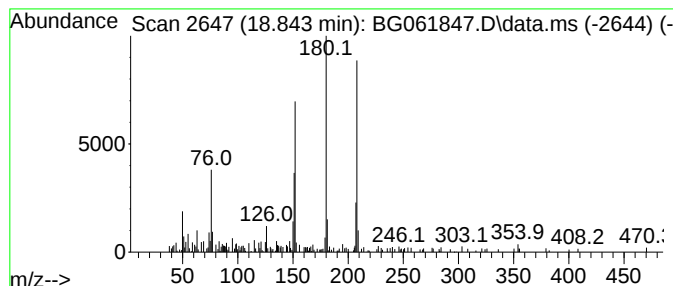
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.843	2.27 ng/ul	116899	Phenanthrene-d10	17.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	89
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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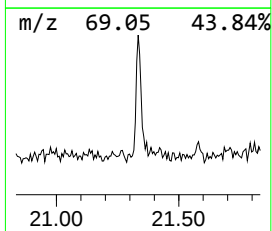
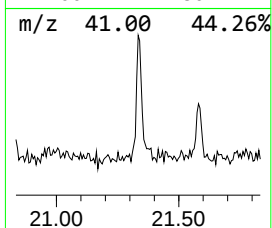
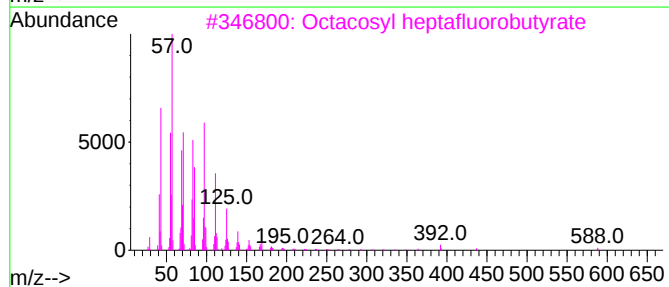
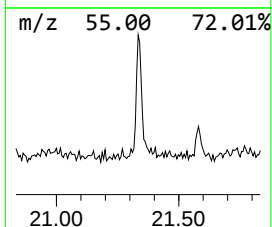
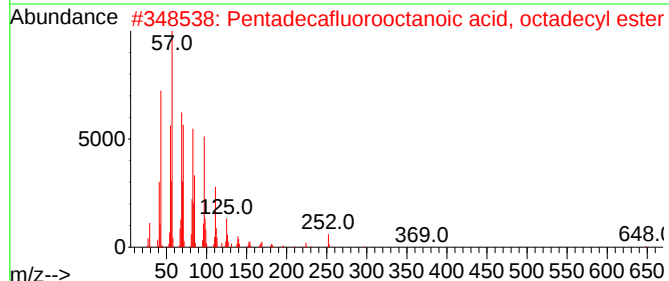
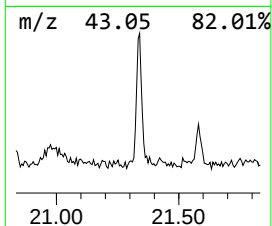
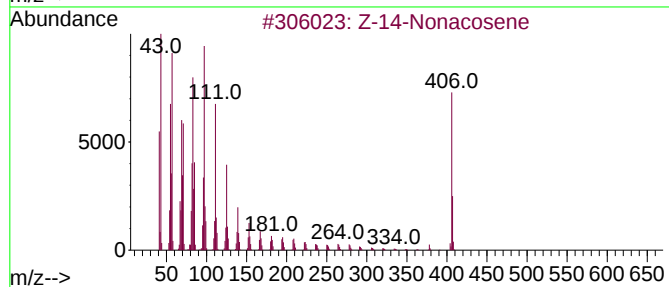
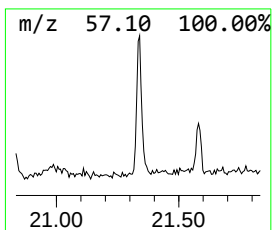
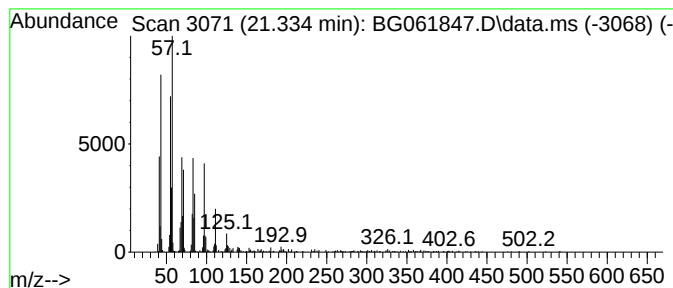
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.334	10.98 ng/ul	779791	Chrysene-d12	21.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-14-Nonacosene	406	C29H58	1010131-18-9	91
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
3			Octacosyl heptafluorobutyrate	606	C32H57F7O2	1010351-83-6	91
4			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
5			Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1010351-81-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070124\
Data File : BG061847.D
Acq On : 2 Jul 2024 15:04
Operator : MA/JU
Sample : P2963-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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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unknown-02	5.135	4.2	ng/ul	98447	1	8.055	473035	20.0
Pentanedioic ac...	9.777	2.4	ng/ul	86011	2	10.870	707467	20.0
1-Phenoxypropan...	11.604	4.1	ng/ul	146505	2	10.870	707467	20.0
Phenanthrene, 2...	18.308	3.3	ng/ul	170069	4	17.433	1027710	20.0
Benzaldehyde, 3...	18.496	4.0	ng/ul	203288	4	17.433	1027710	20.0
9,10-Anthracene...	18.843	2.3	ng/ul	116899	4	17.433	1027710	20.0
(DEL) Alkane: S...	21.334	11.0	ng/ul	779791	5	21.698	1420460	20.0