

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062946.D
 Acq On : 19 Sep 2024 16:49
 Operator : JU/RC
 Sample : P3878-13MEDL2 50X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCVZ1MEDL2

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-BG091724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062946.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.880	303	305	307	rBV3	4122	3417	0.40%	0.065%
2	5.197	357	359	361	rBV2	4416	2686	0.32%	0.051%
3	5.509	408	412	415	rBV5	3145	4564	0.54%	0.087%
4	5.691	441	443	445	rBV	2310	2526	0.30%	0.048%
5	5.820	463	465	467	rBV3	3032	2546	0.30%	0.048%
6	5.885	469	476	483	rBV4	17249	34128	4.02%	0.650%
7	6.055	504	505	508	rVB2	3177	2428	0.29%	0.046%
8	6.084	508	510	512	rVB2	2658	2472	0.29%	0.047%
9	6.267	536	541	544	rBV4	4145	6266	0.74%	0.119%
10	6.349	553	555	564	rBV4	6709	14684	1.73%	0.280%
11	6.419	564	567	571	rBV5	7331	11516	1.36%	0.219%
12	6.472	574	576	577	rBV2	3615	3083	0.36%	0.059%
13	6.748	617	623	631	rBV4	5275	11900	1.40%	0.227%
14	6.837	633	638	639	rBV2	6365	9110	1.07%	0.174%
15	6.913	647	651	654	rBV3	10179	14885	1.75%	0.284%
16	6.948	654	657	661	rVB4	10168	17109	2.02%	0.326%
17	7.025	661	670	675	rBV5	15981	36242	4.27%	0.690%
18	7.083	677	680	686	rVB4	8851	12801	1.51%	0.244%
19	7.136	686	689	691	rBV3	3458	3150	0.37%	0.060%
20	7.248	704	708	712	rBV4	8133	15285	1.80%	0.291%
21	7.483	744	748	757	rVB5	51637	98092	11.56%	1.868%
22	7.724	787	789	792	rBV2	2072	2547	0.30%	0.049%
23	7.853	805	811	818	rBV2	141090	261274	30.78%	4.977%
24	7.918	818	822	828	rVB3	17440	30821	3.63%	0.587%
25	7.965	828	830	833	rBV2	7151	8292	0.98%	0.158%
26	8.205	870	871	875	rVB2	3137	2459	0.29%	0.047%
27	8.288	879	885	886	rBV4	5884	11210	1.32%	0.214%
28	8.499	918	921	922	rBV3	7872	8226	0.97%	0.157%
29	8.623	940	942	946	rVB3	10147	7650	0.90%	0.146%
30	8.963	997	1000	1006	rVB6	6927	8891	1.05%	0.169%
31	9.087	1016	1021	1027	rVB	40848	62556	7.37%	1.192%
32	9.281	1051	1054	1056	rBV4	6160	7037	0.83%	0.134%
33	9.339	1062	1064	1065	rBV2	3510	2532	0.30%	0.048%
34	9.475	1085	1087	1088	rBV	4418	4398	0.52%	0.084%
35	9.639	1113	1115	1116	rBV	3939	2707	0.32%	0.052%
36	9.710	1121	1127	1128	rBV2	3787	4890	0.58%	0.093%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
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37	9.892	1156	1158	1160	rBV	2556	2540	0.30%	0.048%
38	9.980	1169	1173	1174	rBV2	3558	4025	0.47%	0.077%
39	10.009	1174	1178	1181	rVV4	6733	9337	1.10%	0.178%
40	10.068	1187	1188	1193	rBV3	4057	6467	0.76%	0.123%
41	10.344	1232	1235	1238	rBV3	5153	4328	0.51%	0.082%
42	10.421	1244	1248	1249	rBV2	5439	6465	0.76%	0.123%
43	10.520	1263	1265	1268	rBV3	3617	3839	0.45%	0.073%
44	10.644	1276	1286	1293	rBV	185779	374075	44.07%	7.125%
45	10.761	1303	1306	1313	rBV7	9177	14748	1.74%	0.281%
46	10.920	1330	1333	1334	rBV3	4168	3456	0.41%	0.066%
47	11.026	1344	1351	1357	rBV4	21239	36666	4.32%	0.698%
48	11.067	1357	1358	1362	rVB4	3697	4070	0.48%	0.078%
49	11.108	1362	1365	1366	rBV4	2509	2832	0.33%	0.054%
50	11.560	1437	1442	1443	rBV3	6438	8340	0.98%	0.159%
51	11.678	1457	1462	1465	rBV3	9279	16448	1.94%	0.313%
52	11.737	1470	1472	1477	rVV4	10192	15685	1.85%	0.299%
53	11.913	1495	1502	1510	rVB4	12936	24536	2.89%	0.467%
54	12.072	1524	1529	1531	rBV4	13131	20269	2.39%	0.386%
55	12.142	1540	1541	1543	rBV2	2818	2518	0.30%	0.048%
56	12.318	1565	1571	1578	rVB4	25293	45615	5.37%	0.869%
57	12.377	1578	1581	1583	rBV	3361	2771	0.33%	0.053%
58	12.471	1594	1597	1599	rBV2	5964	7593	0.89%	0.145%
59	12.606	1618	1620	1622	rVB2	4522	3653	0.43%	0.070%
60	12.641	1624	1626	1628	rBV2	3149	2806	0.33%	0.053%
61	12.712	1636	1638	1639	rBV	4263	3044	0.36%	0.058%
62	12.882	1665	1667	1668	rBV2	5138	3258	0.38%	0.062%
63	12.965	1676	1681	1682	rBV3	3130	3780	0.45%	0.072%
64	13.241	1724	1728	1730	rBV3	7801	8599	1.01%	0.164%
65	13.323	1735	1742	1748	rBV3	15984	28296	3.33%	0.539%
66	13.405	1754	1756	1759	rBV2	4522	5071	0.60%	0.097%
67	13.446	1761	1763	1768	rVB4	6876	6655	0.78%	0.127%
68	13.523	1774	1776	1777	rBV	3576	2585	0.30%	0.049%
69	13.629	1793	1794	1796	rBV2	3651	3506	0.41%	0.067%
70	13.811	1823	1825	1830	rBV5	5188	7042	0.83%	0.134%
71	13.893	1837	1839	1842	rVB4	10215	8135	0.96%	0.155%
72	13.940	1845	1847	1849	rBV2	4987	5608	0.66%	0.107%
73	14.187	1885	1889	1891	rBV3	7786	8559	1.01%	0.163%
74	14.392	1921	1924	1928	rVB3	26961	33449	3.94%	0.637%
75	14.486	1934	1940	1947	rVB	319811	500539	58.97%	9.534%
76	14.633	1961	1965	1970	rVB3	14676	21966	2.59%	0.418%

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77	14.968	2021	2022	2026	rVB4	7073	5372	0.63%	0.102%
78	15.115	2043	2047	2051	rBV7	24950	39166	4.61%	0.746%
79	15.479	2107	2109	2114	rVB2	10836	14406	1.70%	0.274%
80	15.703	2145	2147	2149	rBV2	3766	3469	0.41%	0.066%
81	15.755	2152	2156	2160	rVV5	11502	15260	1.80%	0.291%
82	16.214	2229	2234	2243	rVB7	17885	44185	5.21%	0.842%
83	16.273	2243	2244	2247	rBV2	4756	3441	0.41%	0.066%
84	16.513	2282	2285	2288	rBV3	5469	8219	0.97%	0.157%
85	16.543	2288	2290	2291	rBV2	6962	6596	0.78%	0.126%
86	16.889	2344	2349	2362	rBV	343149	561096	66.10%	10.688%
87	17.007	2365	2369	2372	rVV5	15123	22444	2.64%	0.428%
88	17.054	2375	2377	2381	rVB3	12899	11488	1.35%	0.219%
89	17.183	2397	2399	2400	rBV2	6657	3428	0.40%	0.065%
90	17.236	2403	2408	2418	rVB	447596	692154	81.54%	13.184%
91	17.336	2422	2425	2428	rVB4	14531	21495	2.53%	0.409%
92	17.365	2428	2430	2431	rBV2	5086	3380	0.40%	0.064%
93	17.530	2454	2458	2459	rBV3	7217	9512	1.12%	0.181%
94	17.741	2491	2494	2499	rVB3	14618	20917	2.46%	0.398%
95	17.918	2520	2524	2528	rBV4	10383	12807	1.51%	0.244%
96	18.441	2608	2613	2617	rVB5	9155	15094	1.78%	0.288%
97	19.563	2802	2804	2807	rVB4	8792	7043	0.83%	0.134%
98	19.645	2812	2818	2823	rBV5	15363	31012	3.65%	0.591%
99	21.490	3126	3132	3143	rBV2	537278	841475	99.13%	16.029%
100	24.528	3642	3649	3661	rVB	314599	848845	100.00%	16.169%

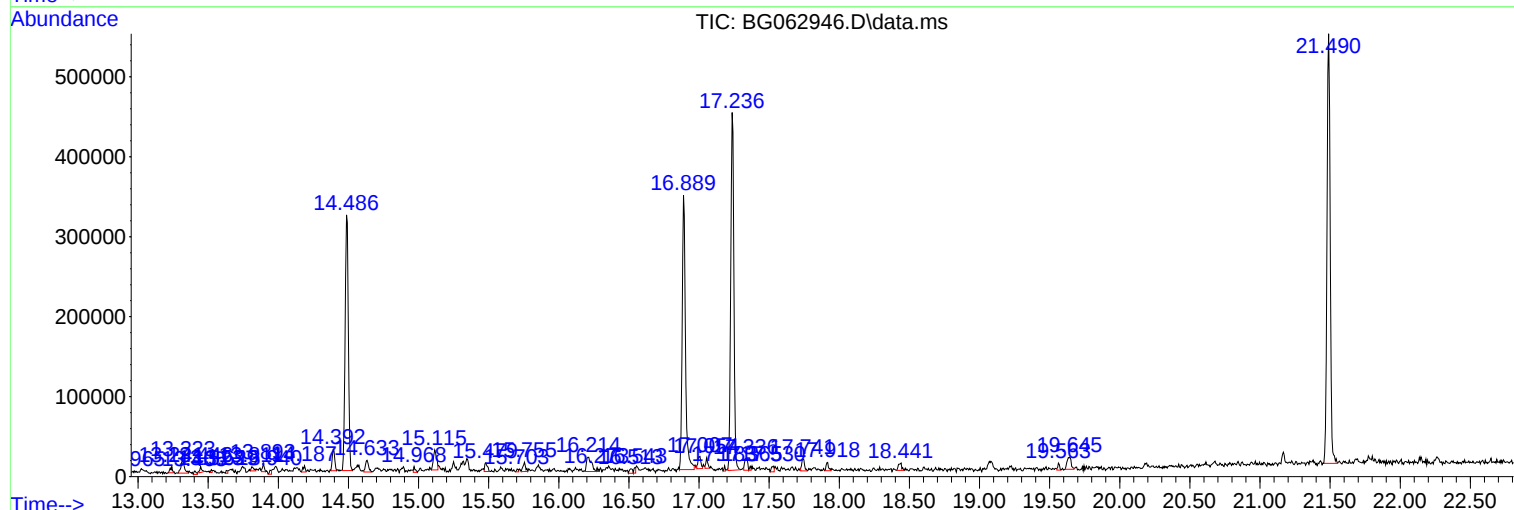
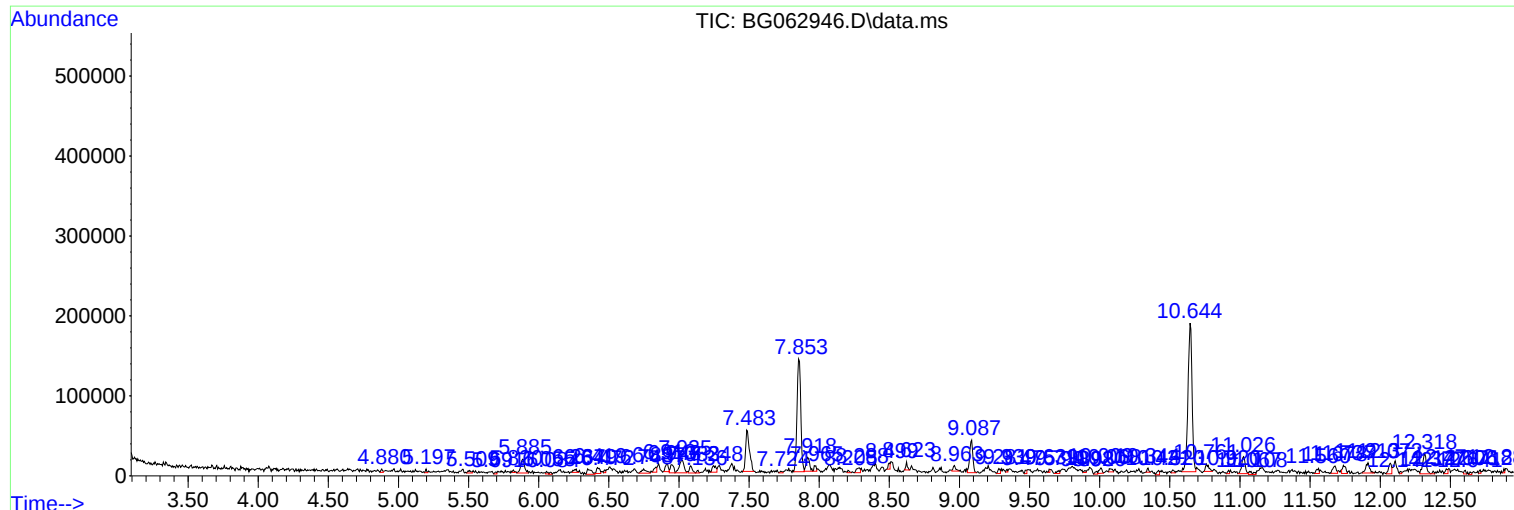
Sum of corrected areas: 5249858

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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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



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Sample : P3878-13MEDL2 50X
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Instrument :
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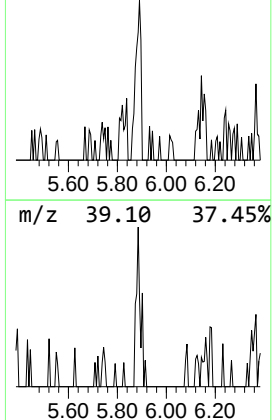
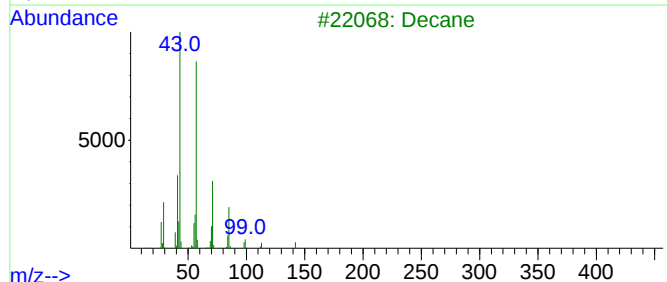
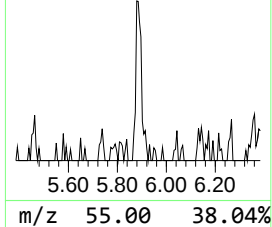
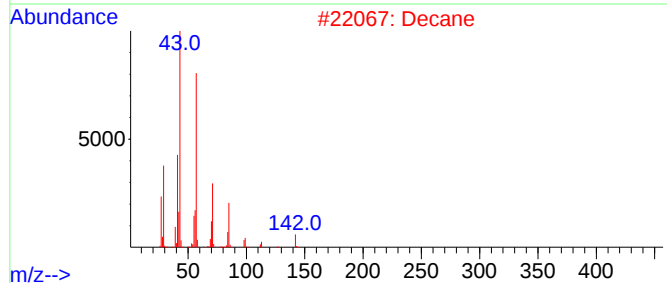
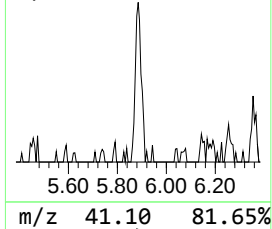
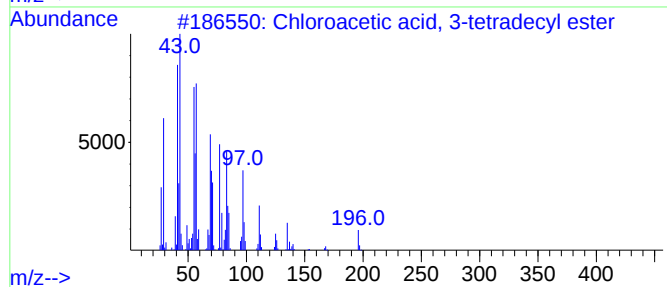
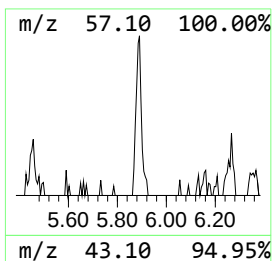
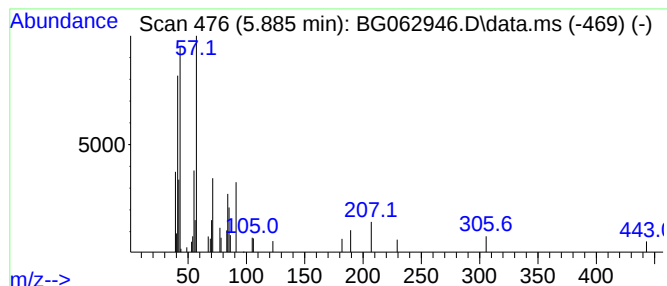
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.885	2.61 ng/ul	34128	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chloroacetic acid, 3-tetradecyl ...	290	C16H31ClO2	1000280-08-5	45
2			Decane	142	C10H22	000124-18-5	43
3			Decane	142	C10H22	000124-18-5	43
4			Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	40
5			Undecane	156	C11H24	001120-21-4	38



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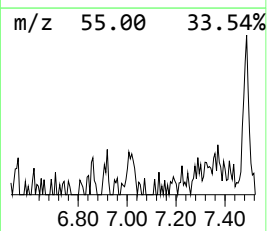
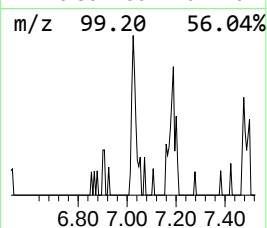
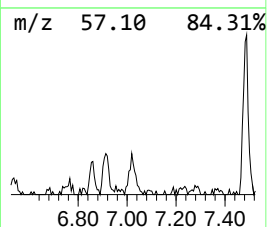
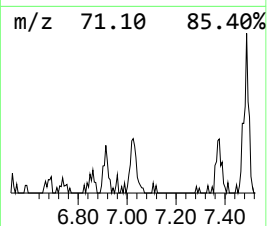
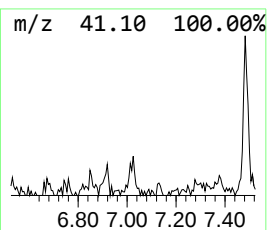
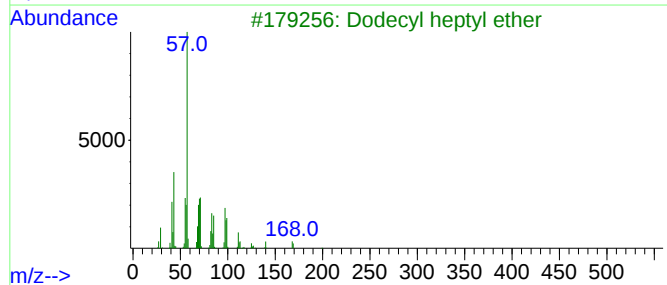
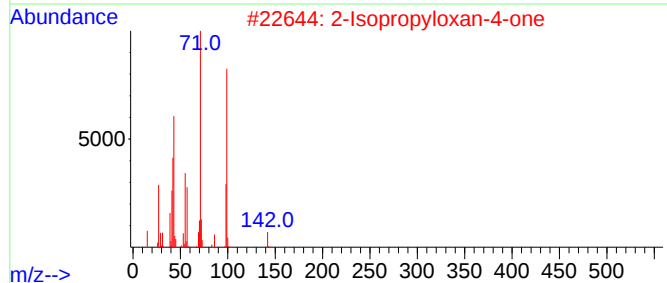
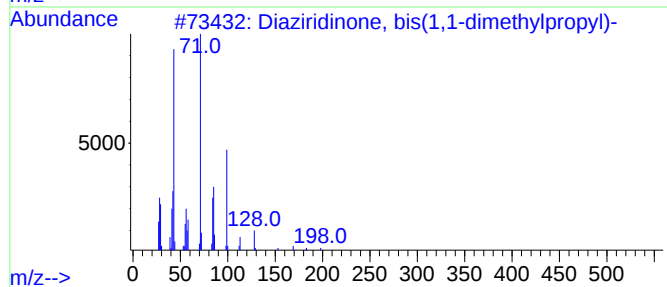
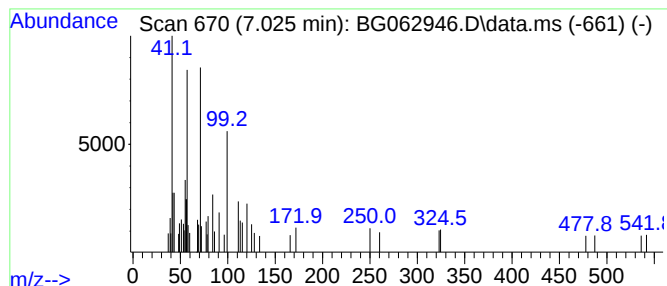
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Diaziridinone, bis(1,1-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.025	2.77 ng/ul	36242	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diaziridinone, bis(1,1-dimethylp...	198	C11H22N2O	019656-75-8	50
2			2-Isopropylloxan-4-one	142	C8H14O2	1000411-56-5	50
3			Dodecyl heptyl ether	284	C19H40O	1010406-39-2	40
4			1-Piperidinyloxy, 4-hydroxy-2,2,...	172	C9H18NO2	002226-96-2	38
5			trans-3-Penten-2-ol	86	C5H10O	003899-34-1	38



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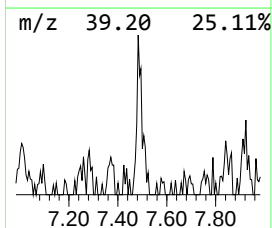
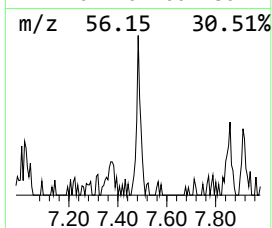
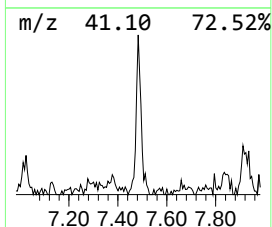
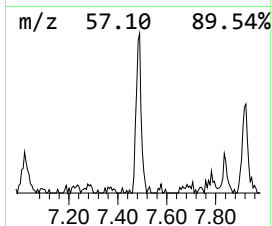
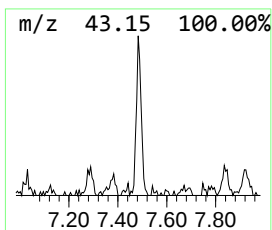
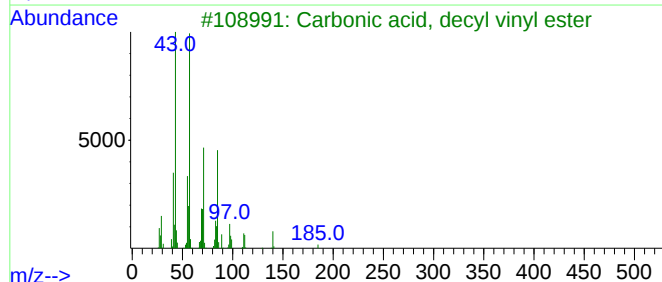
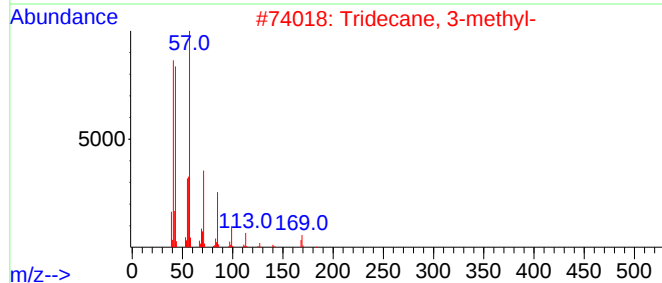
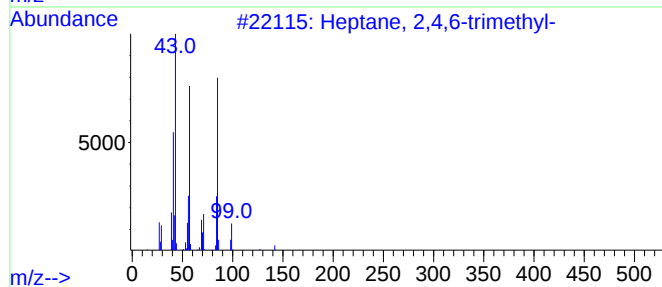
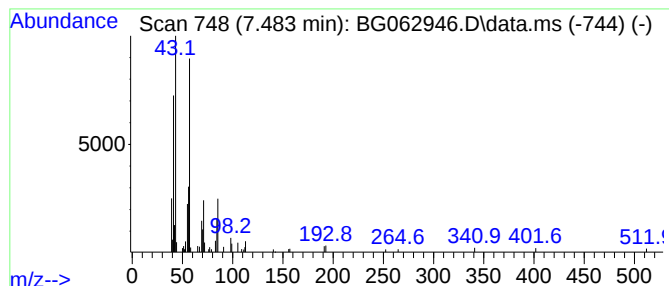
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.483	7.51 ng/ul	98092	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	59
2			Tridecane, 3-methyl-	198	C14H30	006418-41-3	59
3			Carbonic acid, decyl vinyl ester	228	C13H24O3	1000383-25-7	53
4			Undecane	156	C11H24	001120-21-4	53
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062946.D
Acq On : 19 Sep 2024 16:49
Operator : JU/RC
Sample : P3878-13MEDL2 50X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ1MEDL2

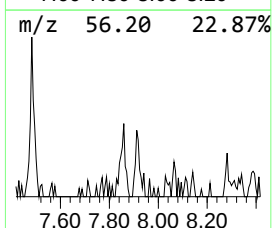
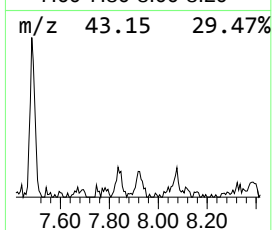
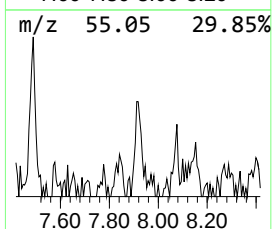
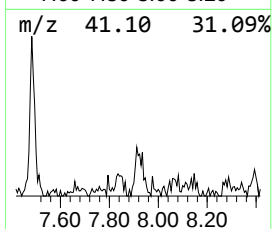
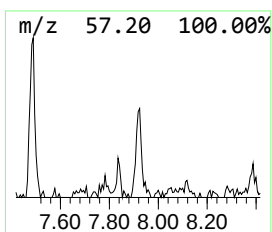
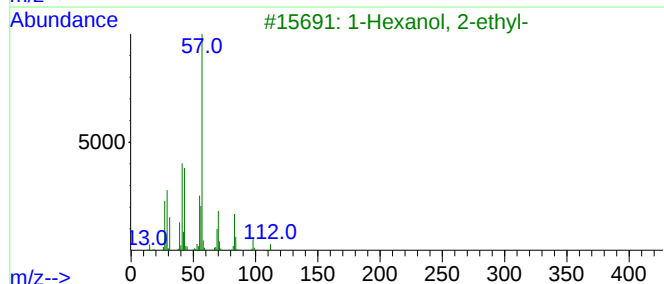
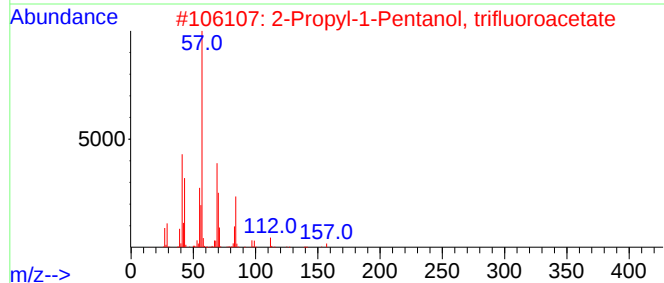
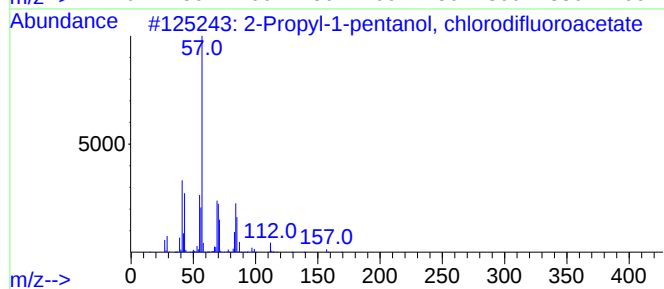
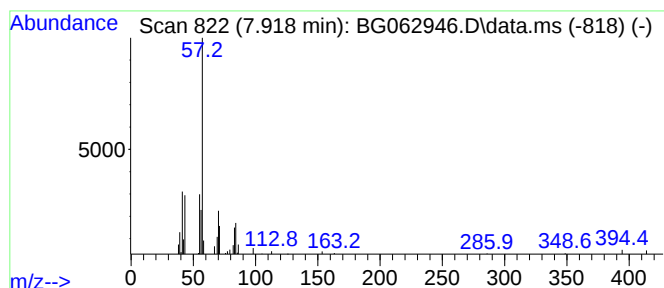
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Propyl-1-pentanol, chloro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.918	2.36 ng/ul	30821	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propyl-1-pentanol, chlorodiflu...	242	C10H17ClF2O2	1000447-27-3	53		
2	2-Propyl-1-Pentanol, trifluoroac...	226	C10H17F3O2	1000365-19-7	53		
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	50		
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	50		
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062946.D
Acq On : 19 Sep 2024 16:49
Operator : JU/RC
Sample : P3878-13MEDL2 50X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ1MEDL2

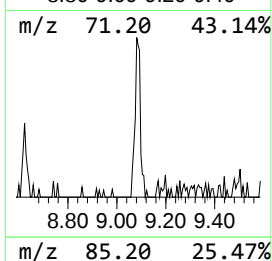
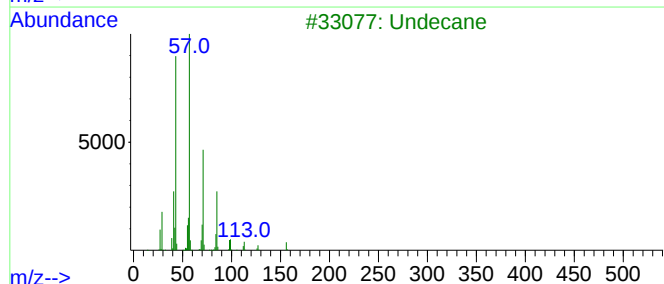
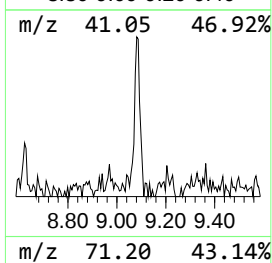
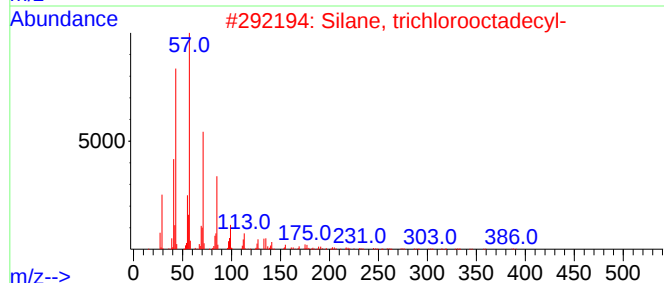
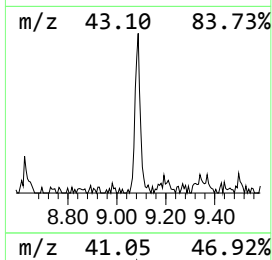
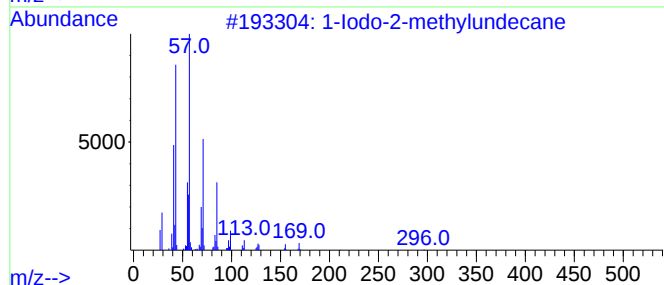
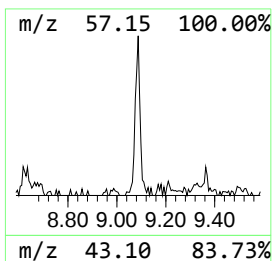
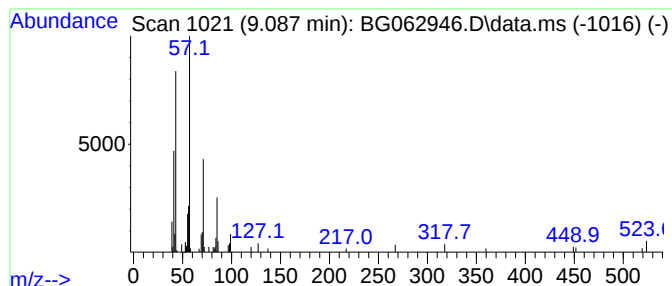
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Iodo-2-methylundecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.087	4.79 ng/ul	62556	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	72
2			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	58
3			Undecane	156	C11H24	001120-21-4	53
4			Tridecane	184	C13H28	000629-50-5	53
5			Pentadecane	212	C15H32	000629-62-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062946.D
Acq On : 19 Sep 2024 16:49
Operator : JU/RC
Sample : P3878-13MEDL2 50X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ1MEDL2

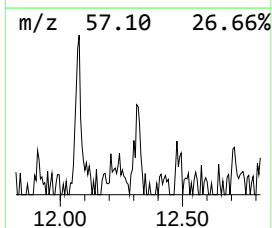
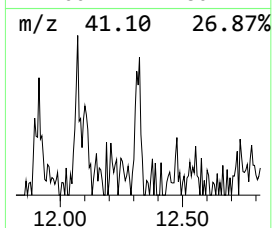
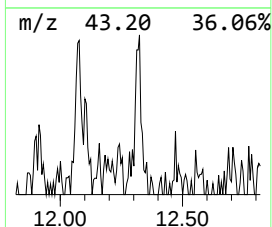
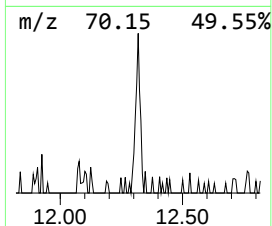
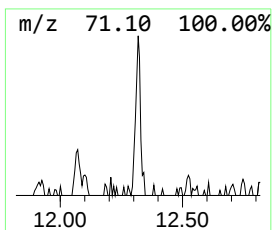
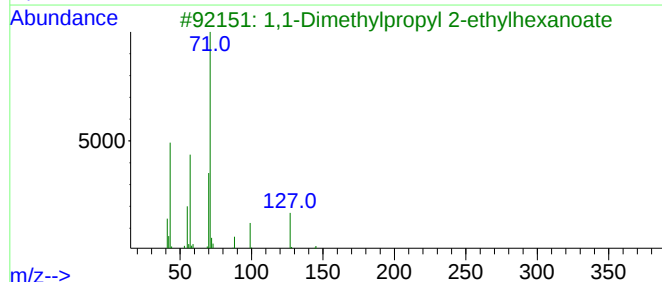
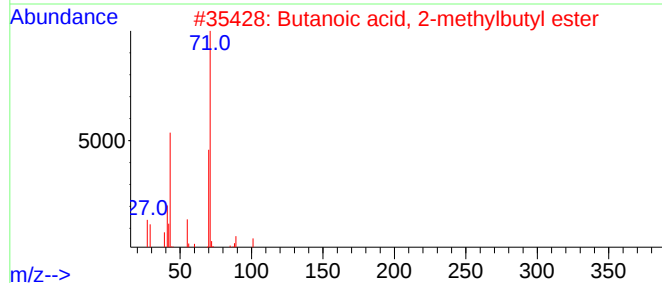
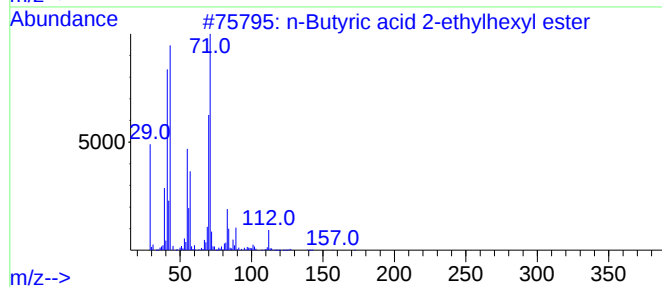
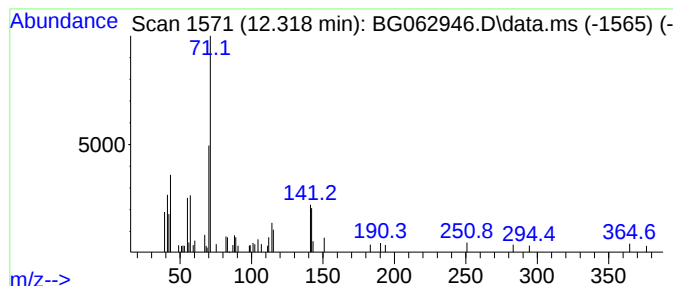
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.318	2.44 ng/ul	45615	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	38
2			Butanoic acid, 2-methylbutyl ester	158	C9H18O2	051115-64-1	38
3			1,1-Dimethylpropyl 2-ethylhexanoate	214	C13H26O2	1000099-99-2	38
4			Octahydro-pyrano[3,2-b]pyran	142	C8H14O2	1000189-97-0	32
5			Butanoic acid, octyl ester	200	C12H24O2	000110-39-4	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062946.D
Acq On : 19 Sep 2024 16:49
Operator : JU/RC
Sample : P3878-13MEDL2 50X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCVZ1MEDL2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Diaziridinone, ...	7.025	2.8	ng/ul	36242	1	7.853	261274	20.0
(DEL) Alkane: S...	7.483	7.5	ng/ul	98092	1	7.853	261274	20.0
2-Propyl-1-pent...	7.918	2.4	ng/ul	30821	1	7.853	261274	20.0
1-Iodo-2-methyl...	9.087	4.8	ng/ul	62556	1	7.853	261274	20.0
unknown-02	12.318	2.4	ng/ul	45615	2	10.644	374075	20.0