

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
 Data File : BG062035.D
 Acq On : 12 Jul 2024 9:22
 Operator : MA/JU
 Sample : P3088-10
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BR8

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: BG062035.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.357	10	12	13	rVB2	5922	4023	0.20%	0.021%
2	3.440	22	26	30	rVB3	10852	15767	0.80%	0.083%
3	3.510	35	38	40	rVB3	4024	3958	0.20%	0.021%
4	3.710	68	72	80	rBV	22900	46965	2.38%	0.247%
5	3.775	82	83	87	rBV4	3871	4276	0.22%	0.023%
6	3.869	96	99	107	rBV2	33080	58017	2.94%	0.305%
7	3.969	112	116	123	rVB	21569	32183	1.63%	0.169%
8	4.092	136	137	142	rVB5	3768	4197	0.21%	0.022%
9	4.204	150	156	158	rBV4	4720	8341	0.42%	0.044%
10	4.756	247	250	255	rBV5	6911	11888	0.60%	0.063%
11	4.956	280	284	288	rBV7	2894	4046	0.21%	0.021%
12	5.114	307	311	316	rVB4	16112	26470	1.34%	0.139%
13	5.208	324	327	333	rBV5	6570	10326	0.52%	0.054%
14	6.025	463	466	469	rVB3	3852	3941	0.20%	0.021%
15	6.084	472	476	477	rBV3	5561	4976	0.25%	0.026%
16	6.577	559	560	564	rVB4	4801	4129	0.21%	0.022%
17	7.036	635	638	641	rBV5	6210	7679	0.39%	0.040%
18	7.141	652	656	657	rBV3	3753	4072	0.21%	0.021%
19	7.200	657	666	677	rVV	242915	423074	21.44%	2.227%
20	7.365	688	694	700	rVB2	136946	222930	11.30%	1.174%
21	7.570	722	729	735	rBV2	217978	381374	19.33%	2.008%
22	7.623	735	738	740	rVV2	10481	8938	0.45%	0.047%
23	8.040	801	809	816	rBV	248271	423896	21.48%	2.232%
24	8.087	816	817	820	rVV3	5724	6500	0.33%	0.034%
25	8.117	820	822	825	rVB4	4703	5354	0.27%	0.028%
26	8.745	922	929	937	rVV2	260056	455996	23.11%	2.400%
27	8.816	939	941	945	rVV4	4345	4633	0.23%	0.024%
28	8.869	945	950	953	rVV5	9935	14554	0.74%	0.077%
29	8.933	958	961	962	rBV2	4629	4537	0.23%	0.024%
30	9.039	976	979	980	rVB3	5094	4087	0.21%	0.022%
31	9.209	1000	1008	1014	rBV	221447	397685	20.15%	2.094%
32	9.356	1031	1033	1035	rBV3	4898	4663	0.24%	0.025%
33	9.580	1068	1071	1074	rBV4	4061	5395	0.27%	0.028%
34	9.679	1085	1088	1090	rBV3	3400	4275	0.22%	0.023%
35	9.756	1096	1101	1107	rVB4	18199	32993	1.67%	0.174%
36	9.926	1123	1130	1137	rBV	208635	378511	19.18%	1.993%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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37	10.473	1216	1223	1230	rBV	338560	574479	29.11%	3.024%
38	10.684	1256	1259	1264	rBV6	3452	5801	0.29%	0.031%
39	10.761	1270	1272	1277	rBV4	3716	4166	0.21%	0.022%
40	10.855	1278	1288	1294	rVV	352653	618597	31.35%	3.256%

41	10.896	1294	1295	1300	rVV4	21478	26480	1.34%	0.139%
42	10.990	1304	1311	1321	rBV2	97394	187801	9.52%	0.989%
43	11.236	1351	1353	1356	rBV3	5114	5832	0.30%	0.031%
44	11.377	1373	1377	1382	rVB4	9839	16176	0.82%	0.085%
45	11.507	1394	1399	1401	rBV3	5499	6307	0.32%	0.033%

46	11.536	1403	1404	1407	rVB2	4615	4207	0.21%	0.022%
47	11.589	1407	1413	1418	rBV2	45439	89129	4.52%	0.469%
48	12.065	1489	1494	1495	rVV3	3112	4064	0.21%	0.021%
49	12.323	1534	1538	1540	rBV3	5038	4972	0.25%	0.026%
50	12.429	1549	1556	1561	rBV6	6409	16845	0.85%	0.089%

51	12.506	1565	1569	1573	rVV7	8777	16302	0.83%	0.086%
52	12.617	1585	1588	1589	rVV	4900	5296	0.27%	0.028%
53	12.641	1590	1592	1596	rVB3	4667	5676	0.29%	0.030%
54	12.717	1603	1605	1613	rVB6	9541	16460	0.83%	0.087%
55	12.934	1641	1642	1646	rVB3	4067	4256	0.22%	0.022%

56	13.111	1669	1672	1675	rBV2	5308	5135	0.26%	0.027%
57	13.422	1722	1725	1727	rBV4	4782	5025	0.25%	0.026%
58	13.481	1733	1735	1738	rVV4	7525	9575	0.49%	0.050%
59	13.528	1741	1743	1744	rVB2	6294	4233	0.21%	0.022%
60	13.587	1750	1753	1754	rBV2	6674	4347	0.22%	0.023%

61	13.810	1787	1791	1792	rBV3	7066	7970	0.40%	0.042%
62	13.992	1817	1822	1825	rBV7	13407	20173	1.02%	0.106%
63	14.074	1828	1836	1842	rBV	604722	836333	42.38%	4.403%
64	14.180	1853	1854	1857	rVV3	5200	5477	0.28%	0.029%
65	14.221	1860	1861	1865	rBV4	5600	7095	0.36%	0.037%

66	14.286	1870	1872	1874	rBV2	3567	4237	0.21%	0.022%
67	14.309	1874	1876	1879	rVV4	7680	8124	0.41%	0.043%
68	14.368	1880	1886	1899	rVB	561639	909776	46.10%	4.789%
69	14.480	1903	1905	1907	rBV2	4069	4191	0.21%	0.022%
70	14.674	1932	1938	1944	rVB	529423	809940	41.05%	4.264%

71	14.732	1944	1948	1955	rVB4	25608	46017	2.33%	0.242%
72	14.791	1955	1958	1965	rBV6	10746	12991	0.66%	0.068%
73	14.879	1965	1973	1984	rBV	304426	563996	28.58%	2.969%
74	14.962	1984	1987	1989	rBV4	6490	7921	0.40%	0.042%
75	14.985	1989	1991	1993	rVB3	6338	4635	0.23%	0.024%

76	15.020	1993	1997	2002	rBV6	24925	32965	1.67%	0.174%
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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	15.067	2002	2005	2011	rVV4	23827	37143	1.88%	0.196%
78	15.285	2038	2042	2044	rBV5	6873	8652	0.44%	0.046%
79	15.667	2100	2107	2112	rBV	722382	1129858	57.26%	5.948%
80	15.719	2112	2116	2120	rVV2	31795	51413	2.61%	0.271%
81	15.778	2121	2126	2133	rVB	182298	264282	13.39%	1.391%
82	15.884	2142	2144	2145	rBV2	8415	6212	0.31%	0.033%
83	15.919	2147	2150	2156	rVB5	30485	45135	2.29%	0.238%
84	16.360	2223	2225	2227	rBV3	12438	14204	0.72%	0.075%
85	16.642	2270	2273	2277	rBV4	22889	32480	1.65%	0.171%
86	17.071	2342	2346	2352	rVB3	29501	44126	2.24%	0.232%
87	17.153	2355	2360	2366	rBV9	28641	58569	2.97%	0.308%
88	17.235	2372	2374	2377	rVB	26115	24419	1.24%	0.129%
89	17.417	2399	2405	2409	rBV	616918	993890	50.37%	5.232%
90	17.459	2409	2412	2417	rVV	457434	642902	32.58%	3.384%
91	17.517	2417	2422	2432	rVB	808498	1282365	64.99%	6.751%
92	18.293	2550	2554	2558	rBV2	183250	249284	12.63%	1.312%
93	18.340	2559	2562	2566	rVV	70745	100190	5.08%	0.527%
94	18.481	2582	2586	2590	rBV4	100554	174238	8.83%	0.917%
95	18.828	2642	2645	2650	rVB2	115710	145221	7.36%	0.764%
96	19.468	2749	2754	2759	rVB	1050661	1429932	72.46%	7.528%
97	19.797	2805	2810	2813	rBV2	713144	1161355	58.85%	6.114%
98	19.826	2813	2815	2823	rVB	536464	698237	35.38%	3.676%
99	21.554	3106	3109	3113	rVB	384690	470945	23.87%	2.479%
100	21.671	3123	3129	3135	rBV3	859544	1973282	100.00%	10.388%

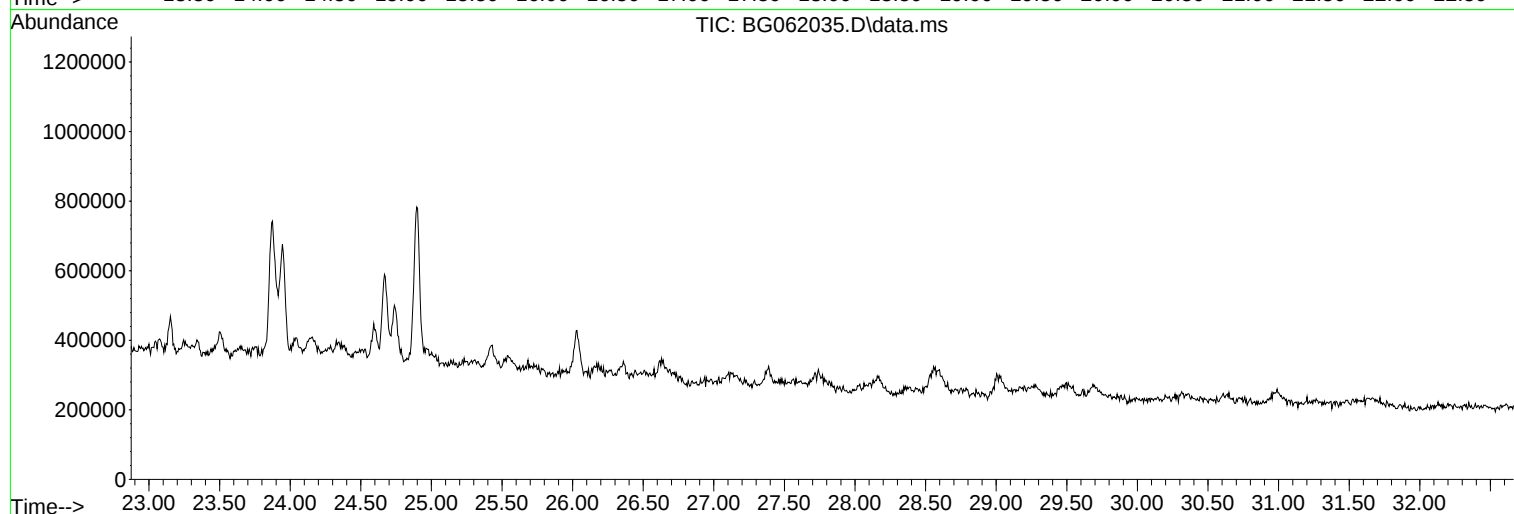
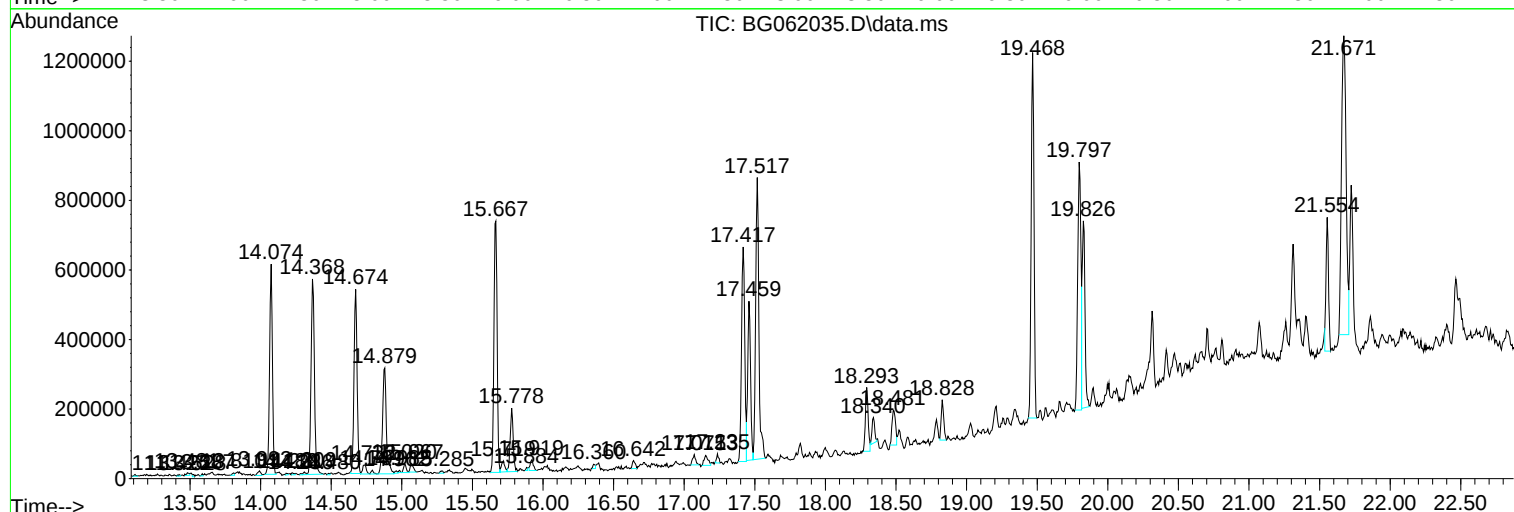
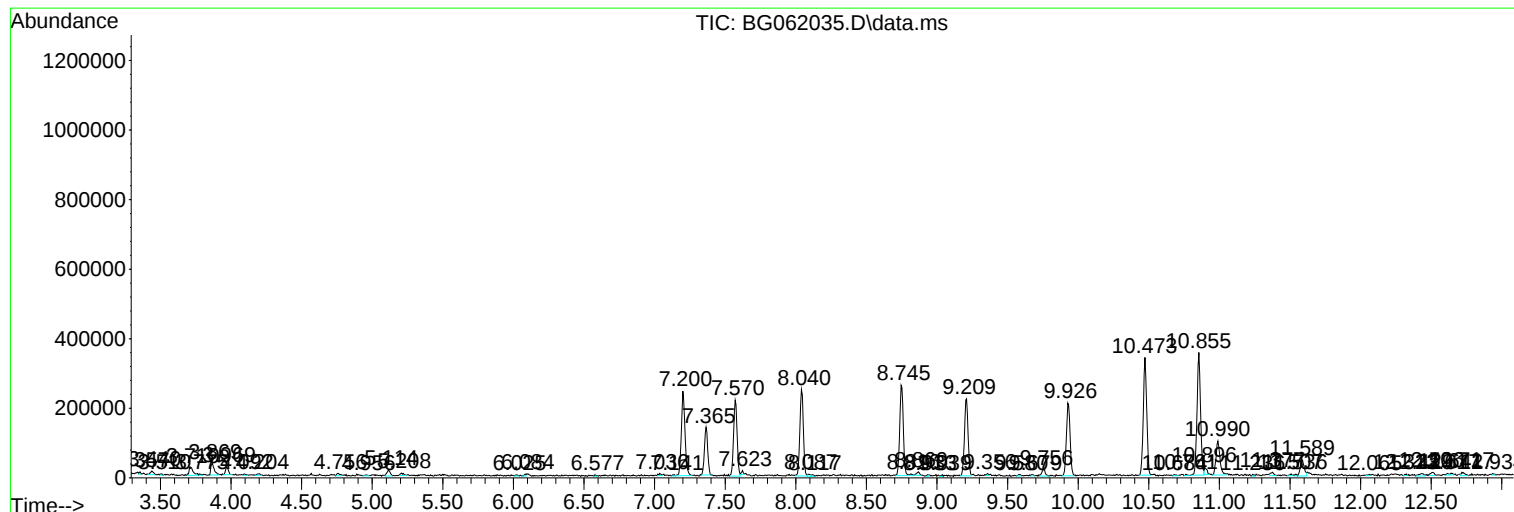
Sum of corrected areas: 18995985

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Misc :
ALS Vial : 32 Sample Multiplier: 1

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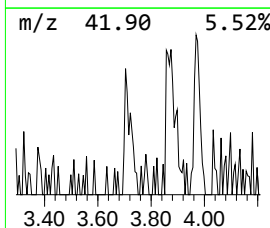
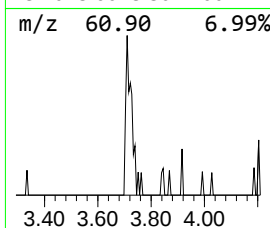
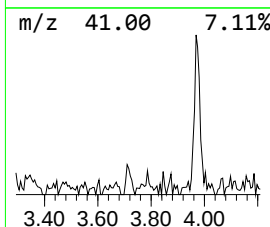
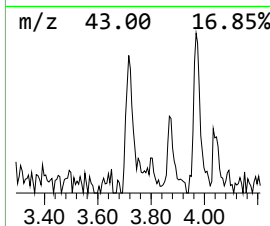
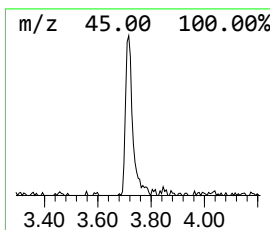
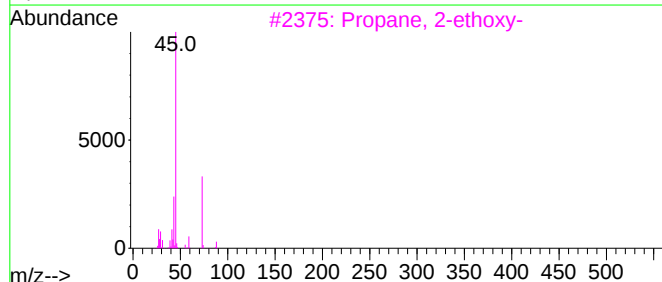
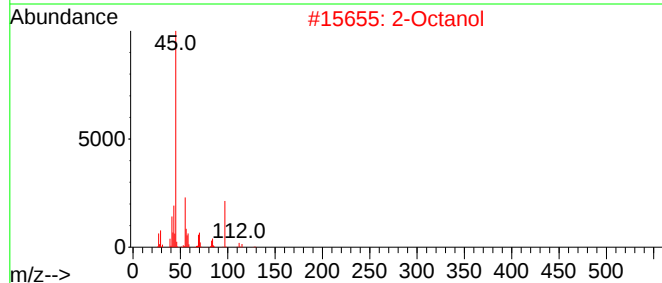
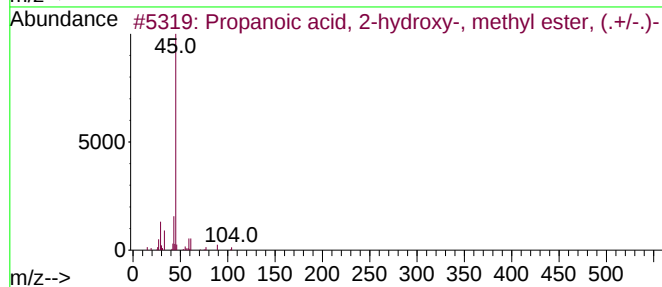
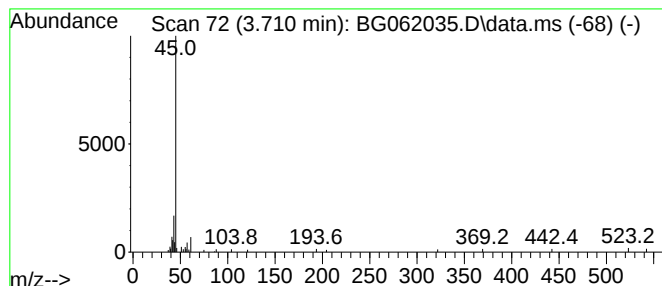
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.710	2.22 ng/ul	46965	1,4-Dichlorobenzene-d4	8.040

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	39
2			2-Octanol	130	C8H18O	000123-96-6	38
3			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	38
4			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	38
5			Butane, 1-methoxy-	88	C5H12O	000628-28-4	9



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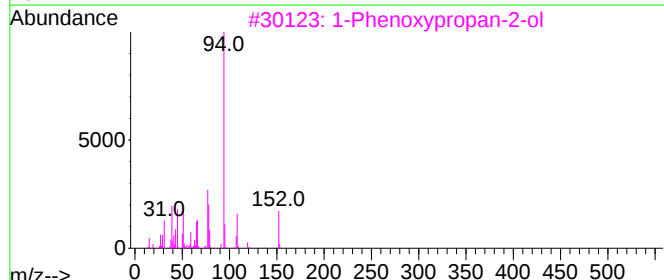
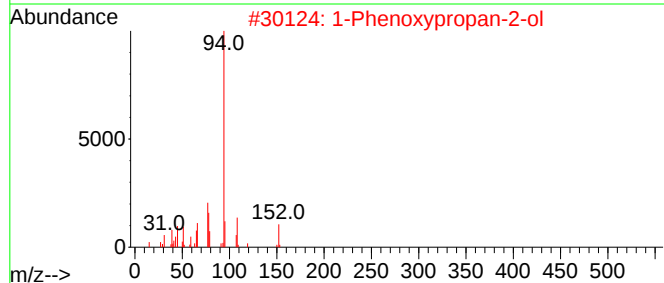
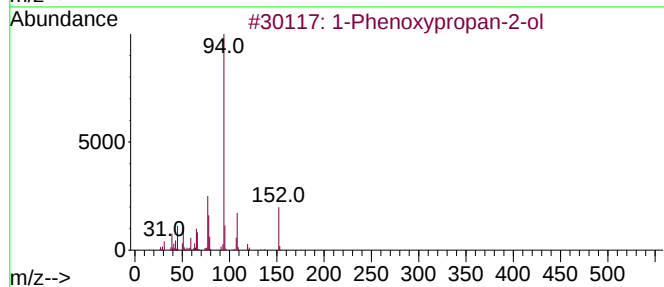
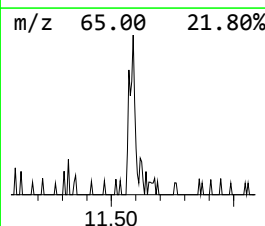
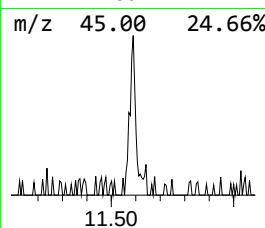
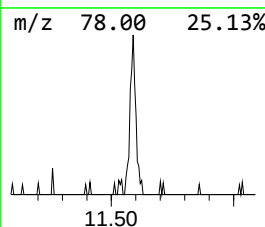
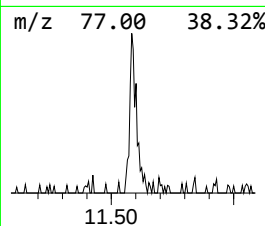
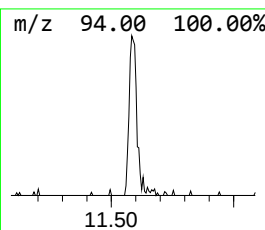
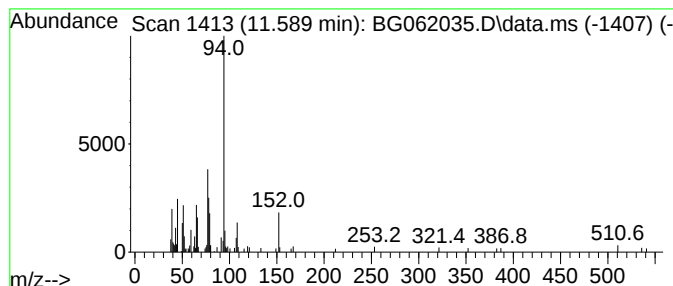
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.589	2.88 ng/ul	89129	Naphthalene-d8	10.855

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	89
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	76
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	58
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	52



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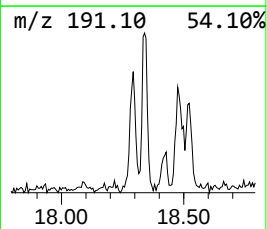
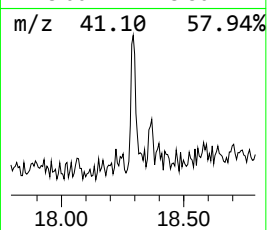
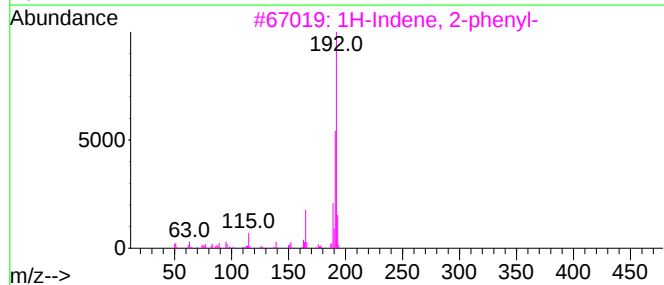
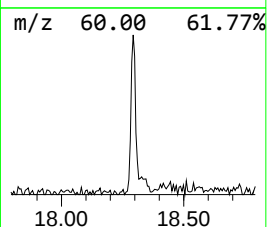
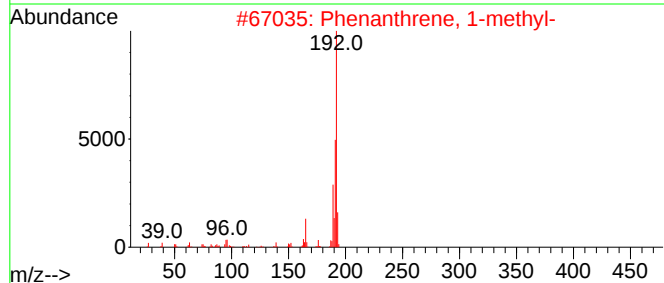
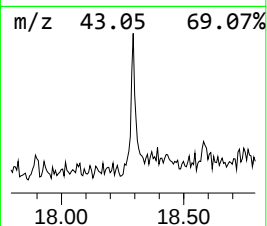
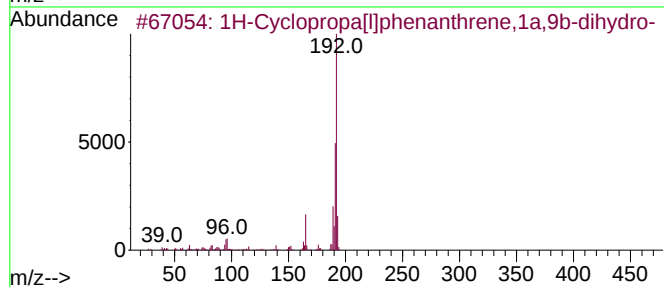
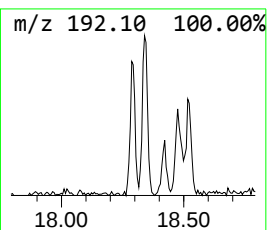
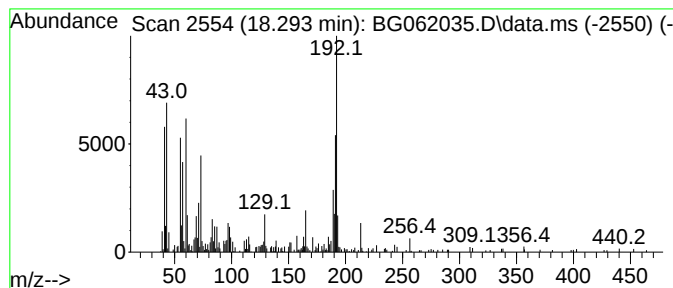
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1H-Cyclopropa[l]phenanthren... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.293	5.02 ng/ul	249284	Phenanthrene-d10	17.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	92
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	83
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062035.D
Acq On : 12 Jul 2024 9:22
Operator : MA/JU
Sample : P3088-10
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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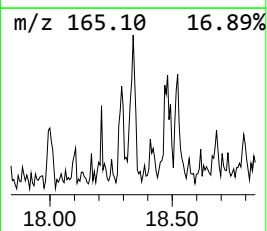
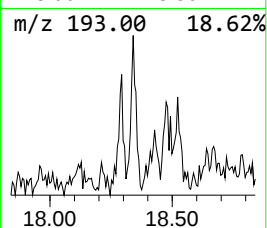
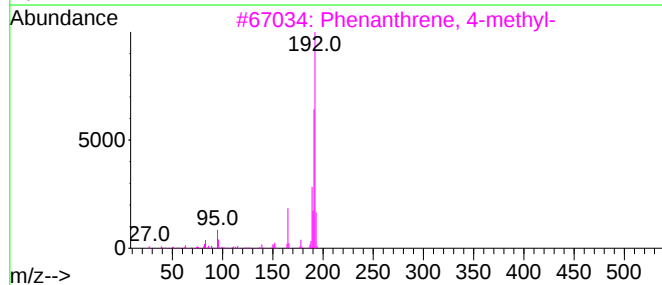
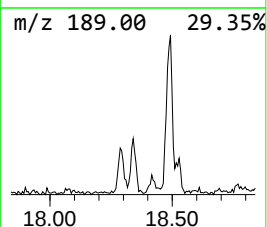
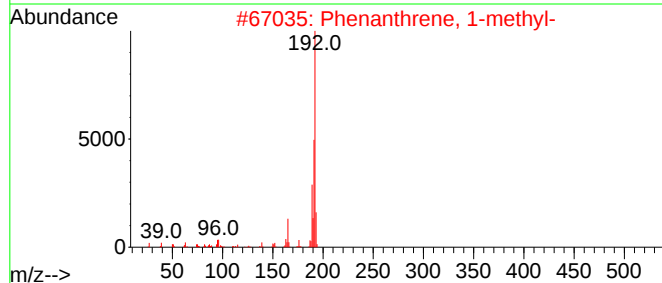
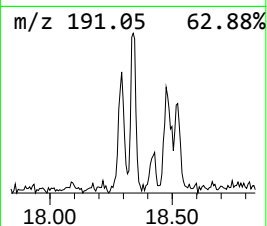
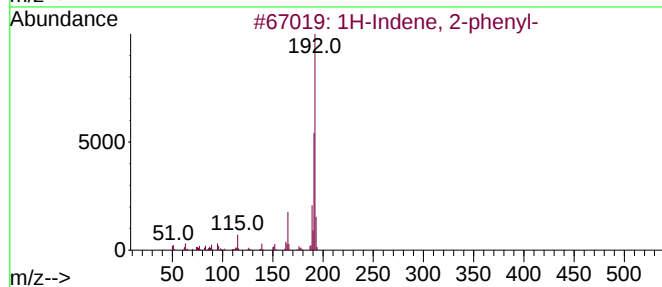
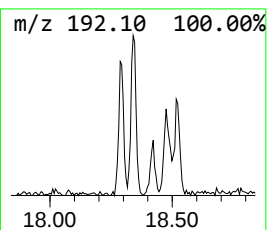
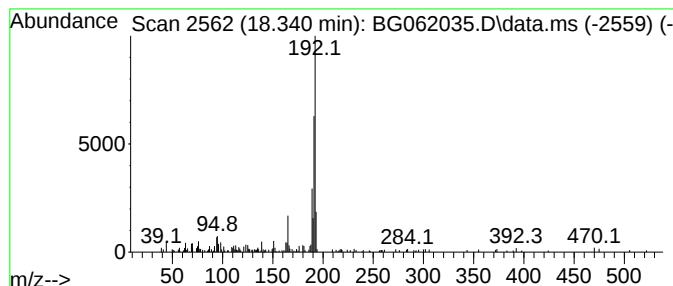
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1H-Indene, 2-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.340	2.02 ng/ul	100190	Phenanthrene-d10	17.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-			192	C15H12	004505-48-0	95
2	Phenanthrene, 1-methyl-			192	C15H12	000832-69-9	95
3	Phenanthrene, 4-methyl-			192	C15H12	000832-64-4	94
4	1H-Indene, 2-phenyl-			192	C15H12	004505-48-0	94
5	1H-Cyclopropa[1]phenanthrene,1a,...			192	C15H12	000949-41-7	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062035.D
Acq On : 12 Jul 2024 9:22
Operator : MA/JU
Sample : P3088-10
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BR8

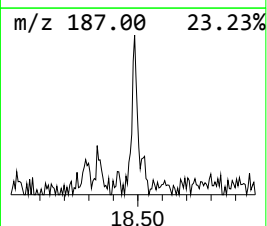
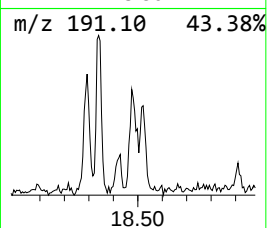
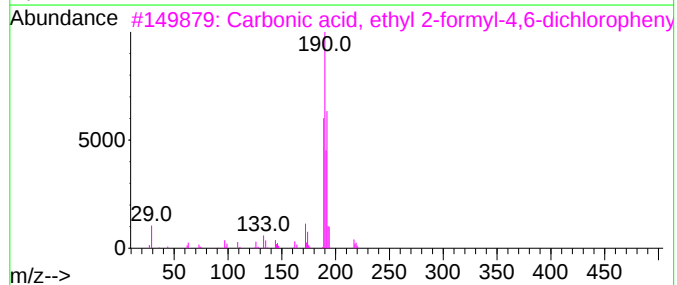
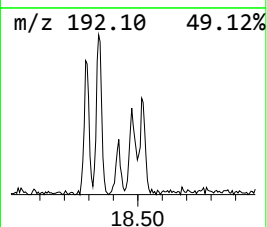
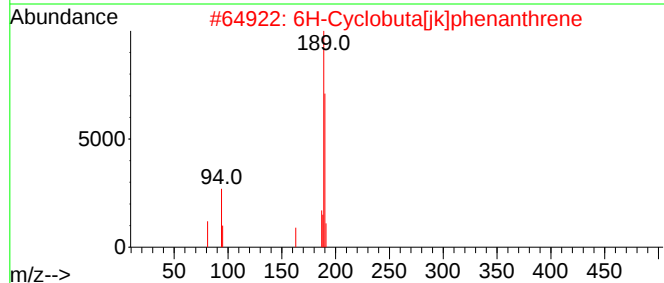
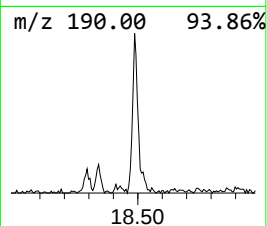
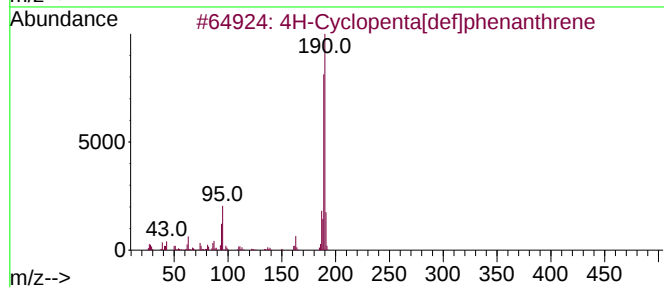
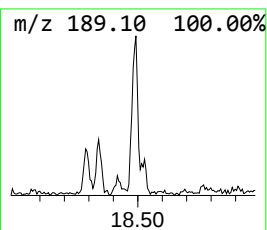
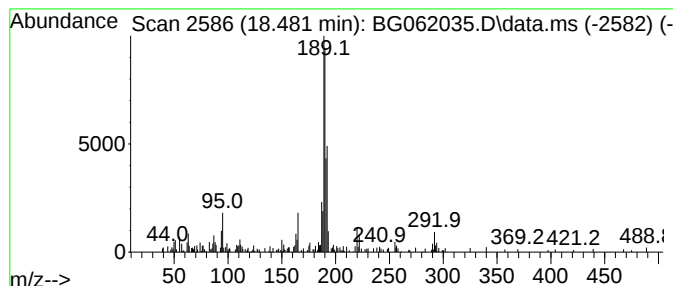
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.481	3.51 ng/ul	174238	Phenanthrene-d10	17.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
3			Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	40
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	40
5			Carbonic acid, 2-formyl-4,6-dich...	276	C11H10Cl2O4	1000331-39-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062035.D
Acq On : 12 Jul 2024 9:22
Operator : MA/JU
Sample : P3088-10
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BR8

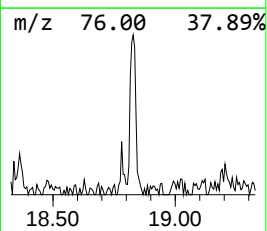
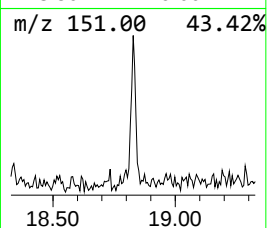
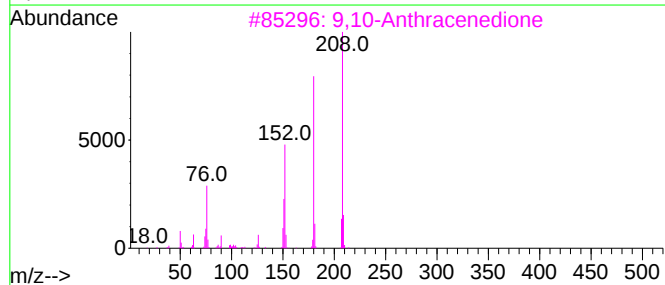
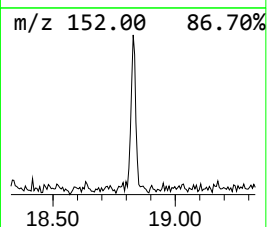
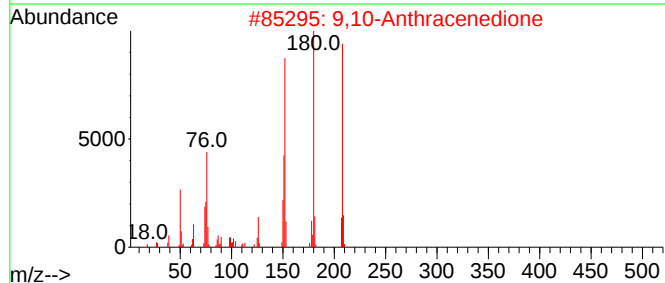
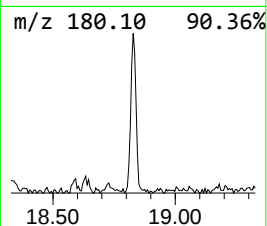
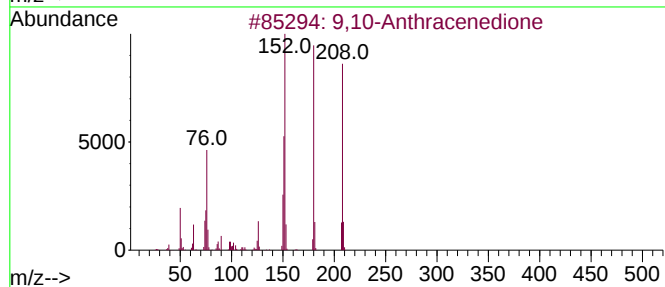
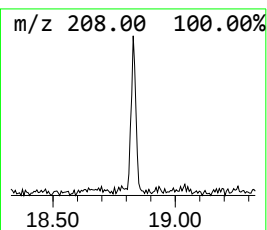
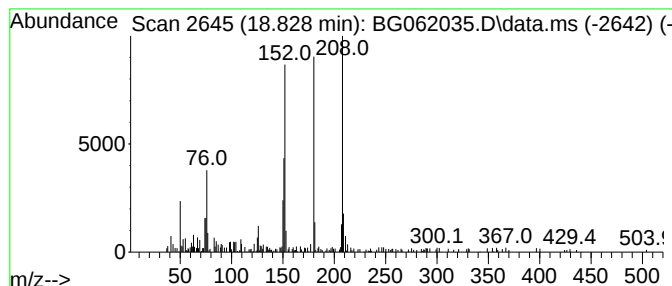
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.828	2.92 ng/ul	145221	Phenanthrene-d10	17.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	91
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	91
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062035.D
Acq On : 12 Jul 2024 9:22
Operator : MA/JU
Sample : P3088-10
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

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1-Phenoxypropan...	11.589	2.9	ng/u1	89129	2	10.855	618597	20.0
1H-Cyclopropa[1...	18.293	5.0	ng/u1	249284	4	17.417	993890	20.0
1H-Indene, 2-ph...	18.340	2.0	ng/u1	100190	4	17.417	993890	20.0
4H-Cyclopenta[d...	18.481	3.5	ng/u1	174238	4	17.417	993890	20.0
9,10-Anthracene...	18.828	2.9	ng/u1	145221	4	17.417	993890	20.0