LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

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-BG071924.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062200.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.883	97	101	108	rBV	58659	97491	4.58%	0.463%
2	5.363	347	353	363	rVV	223333	374232	17.60%	1.776%
3	6.538	547	553	562	rVB2	28166	55535	2.61%	0.264%
4	7.031	630	637	643	rVV2	36993	66550	3.13%	0.316%
5	7.196	660	665	669	rVV5	21824	37162	1.75%	0.176%
6	7.249	669	674	684	rVB2	80734	149870	7.05%	0.711%
7	7.390	691	698	704	rVV	112151	191807	9.02%	0.910%
8	7.443	704	707	718	rVB3	14641	25770	1.21%	0.122%
9	7.607	728	735	742	rVV2	173905	297384	13.98%	1.411%
10	7.707	746	752	757	rVV2	60243	98906	4.65%	0.469%
11	7.954	784	794	801	rBV4	26141	61658	2.90%	0.293%
12	8.065	806	813	827	rVV2	163712	329102	15.48%	1.562%
13	8.177	827	832	840	rVV3	24423	48069	2.26%	0.228%
14	8.365	857	864	868	rBV2	8902	14256	0.67%	0.068%
15	8.435	868	876	884	rVB2	40193	79310	3.73%	0.376%
16	8.547	891	895	900	rVB3	18076	30312	1.43%	0.144%
17	8.700	915	921	927	rBV3	23396	43416	2.04%	0.206%
18	8.794	931	937	945	rVV	206251	353293	16.61%	1.676%
19	9.011	969	974	980	rBV6	10645	20775	0.98%	0.099%
20	9.170	994	1001	1006	rBV3	90322	157731	7.42%	0.748%
21	9.240	1006	1013	1022	rVB2	245226	479162	22.53%	2.274%
22	9.411	1035	1042	1048	rBV5	11539	24875	1.17%	0.118%
23	9.510	1053	1059	1067	rVB2	25916	52113	2.45%	0.247%
24	9.692	1084	1090	1096	rVB2	42136	68914	3.24%	0.327%
25	9.751	1096	1100	1104	rBV3	20125	37138	1.75%	0.176%
26	9.957	1128	1135	1143	rBV2	168682	308709	14.52%	1.465%
27	10.057	1145	1152	1158	rBV5	26375	48475	2.28%	0.230%
28	10.115	1158	1162	1166	rBV2	18930	30051	1.41%	0.143%
29	10.268	1180	1188	1195	rVB3	117749	214628	10.09%	1.018%
30	10.444	1210	1218	1222	rBV7	13775	33570	1.58%	0.159%
31	10.515	1222	1230	1237	rBV2	286209	509328	23.95%	2.417%
32	10.809	1271	1280	1282	rBV6	10678	27082	1.27%	0.129%
33	10.885	1282	1293	1298	rBV	227704	480118	22.58%	2.278%
34	10.985	1305	1310	1311	rBV3	30474	44414	2.09%	0.211%
35	11.020	1311	1316	1323	rVV	164026	297230	13.98%	1.410%
36	11.067	1323	1324	1330	rVB4	11158	16881	0.79%	0.080%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

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-BG071924.MA.M
Title : SVOA CALIBRATION

37	11.455	1385	1390	1394	rVV6	10850	18793	0.88%	0.089%
38	11.531	1399	1403	1409	rBV5	13161	26814	1.26%	0.127%
39	11.619	1412	1418	1423	rVV4	13505	28050	1.32%	0.133%
40	11.778	1441	1445	1451	rVB2	30037	53078	2.50%	0.252%
41	11.978	1473	1479	1483	rVV2	27332	46537	2.19%	0.221%
42	12.054	1487	1492	1501	rVB7	16519	37450	1.76%	0.178%
43	12.189	1510	1515	1520	rVB2	17886	25446	1.20%	0.121%
44	12.254	1521	1526	1532	rVB3	29968	48750	2.29%	0.231%
45	12.412	1548	1553	1560	rVV8	32000	57981	2.73%	0.275%
46	12.530	1567	1573	1580	rVV	542113	903872	42.50%	4.289%
47	12.594	1580	1584	1589	rVB6	18518	33227	1.56%	0.158%
48	12.753	1601	1611	1618	rVV	1243087	2126621	100.00%	10.091%
49	12.947	1640	1644	1648	rVV7	12895	21167	1.00%	0.100%
50	13.047	1654	1661	1663	rVB8	13648	28480	1.34%	0.135%
51	13.288	1698	1702	1706	rVB7	8411	15297	0.72%	0.073%
52	13.493	1732	1737	1745	rBV	65085	101721	4.78%	0.483%
53	13.622	1755	1759	1765	rBV3	31953	52789	2.48%	0.250%
54	13.722	1772	1776	1777	rBV4	33512	31572	1.48%	0.150%
55	13.857	1789	1799	1801	rBV	78304	146143	6.87%	0.693%
56	13.922	1808	1810	1816	rVB7	9050	15351	0.72%	0.073%
57	14.016	1820	1826	1831	rVV3	169931	334295	15.72%	1.586%
58	14.098	1831	1840	1846	rVB	490800	810332	38.10%	3.845%
59	14.251	1861	1866	1870	rBV	76858	135107	6.35%	0.641%
60	14.286	1870	1872	1876	rVB3	36375	41212	1.94%	0.196%
61	14.392	1884	1890	1903	rVB2	463861	851781	40.05%	4.042%
62	14.562	1914	1919	1923	rBV2	38385	53470	2.51%	0.254%
63	14.697	1931	1942	1948	rBV	431112	705243	33.16%	3.347%
64	14.756	1948	1952	1958	rVV2	28688	53222	2.50%	0.253%
65	14.868	1966	1971	1975	rVV7	11307	24097	1.13%	0.114%
66	14.932	1977	1982	1987	rVV3	51778	91441	4.30%	0.434%
67	15.009	1991	1995	1997	rVV5	18549	28400	1.34%	0.135%
68	15.044	1999	2001	2006	rVV6	16072	31697	1.49%	0.150%
69	15.091	2007	2009	2016	rVB8	18893	31787	1.49%	0.151%
70	15.162	2016	2021	2024	rBV7	8949	16350	0.77%	0.078%
71	15.308	2043	2046	2050	rVV6	18028	28710	1.35%	0.136%
72	15.467	2069	2073	2076	rVV5	9752	16917	0.80%	0.080%
73	15.514	2076	2081	2088	rVB6	30355	51890	2.44%	0.246%
74	15.684	2104	2110	2116	rVB	683926	1003175	47.17%	4.760%
75	15.743	2117	2120	2125	rVB6	12457	16480	0.77%	0.078%
76	15.814	2127	2132	2136	rBV2	79858	135262	6.36%	0.642%

LSC Area Percent Report

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Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

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-BG071924.MA.M
Title : SVOA CALIBRATION

77	16.072	2170	2176	2182	rBV4	100780	215648	10.14%	1.023%
78	16.419	2230	2235	2241	rVB2	44472	72182	3.39%	0.343%
79	16.513	2246	2251	2255	rBV	65604	101055	4.75%	0.480%
80	16.566	2255	2260	2266	rVV2	106010	202069	9.50%	0.959%
81	16.630	2266	2271	2275	rVV2	91306	143537	6.75%	0.681%
82	16.671	2275	2278	2282	rVV	50696	68643	3.23%	0.326%
83	16.724	2282	2287	2289	rVV5	21785	35575	1.67%	0.169%
84	16.754	2289	2292	2294	rVV	28699	41499	1.95%	0.197%
85	16.777	2294	2296	2299	rVV4	25150	30020	1.41%	0.142%
86	16.830	2300	2305	2312	rVV3	83310	153551	7.22%	0.729%
87	16.894	2312	2316	2319	rVV2	60973	82437	3.88%	0.391%
88	16.936	2319	2323	2326	rVV6	31638	54724	2.57%	0.260%
89	16.971	2326	2329	2333	rVB2	40833	54622	2.57%	0.259%
90	17.235	2367	2374	2381	rVV2	33075	64924	3.05%	0.308%
91	17.441	2403	2409	2415	rBV	538454	810060	38.09%	3.844%
92	17.541	2420	2426	2433	rVB	723244	1093419	51.42%	5.189%
93	18.304	2554	2556	2563	rVB6	14020	23282	1.09%	0.110%
94	18.569	2598	2601	2607	rBV2	29344	40427	1.90%	0.192%
95	18.992	2670	2673	2678	rVB7	16975	24539	1.15%	0.116%
96	19.820	2809	2814	2821	rVB	911356	1298598	61.06%	6.162%
97	21.700	3129	3134	3141	rVB2	497082	844061	39.69%	4.005%
98	23.350	3410	3415	3422	rVB3	61270	129915	6.11%	0.616%
99	24.713	3639	3647	3658	rVB	423297	1140635	53.64%	5.413%
100	24.937	3677	3685	3696	rVB2	305345	863126	40.59%	4.096%

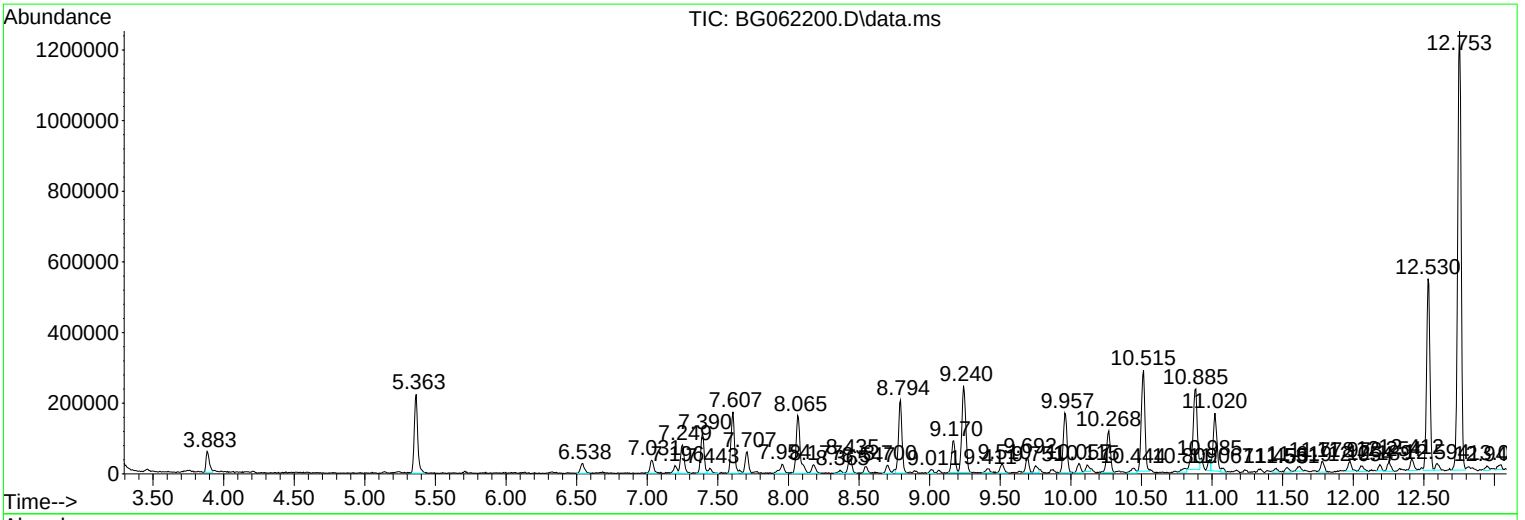
Sum of corrected areas: 21073872

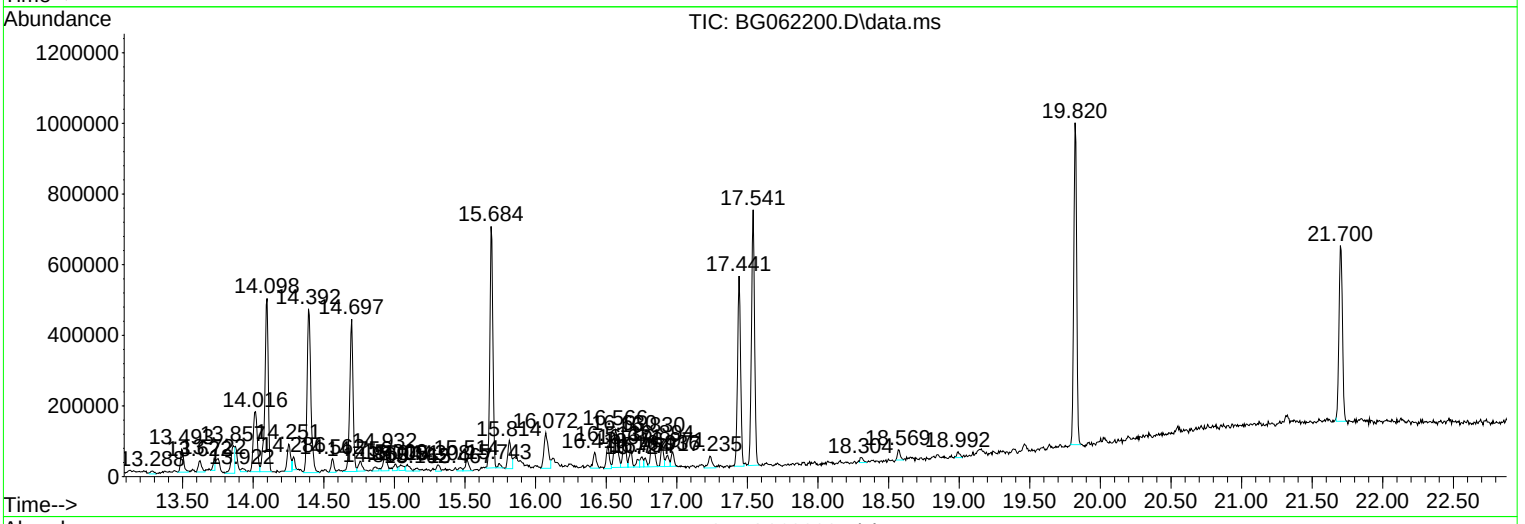
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

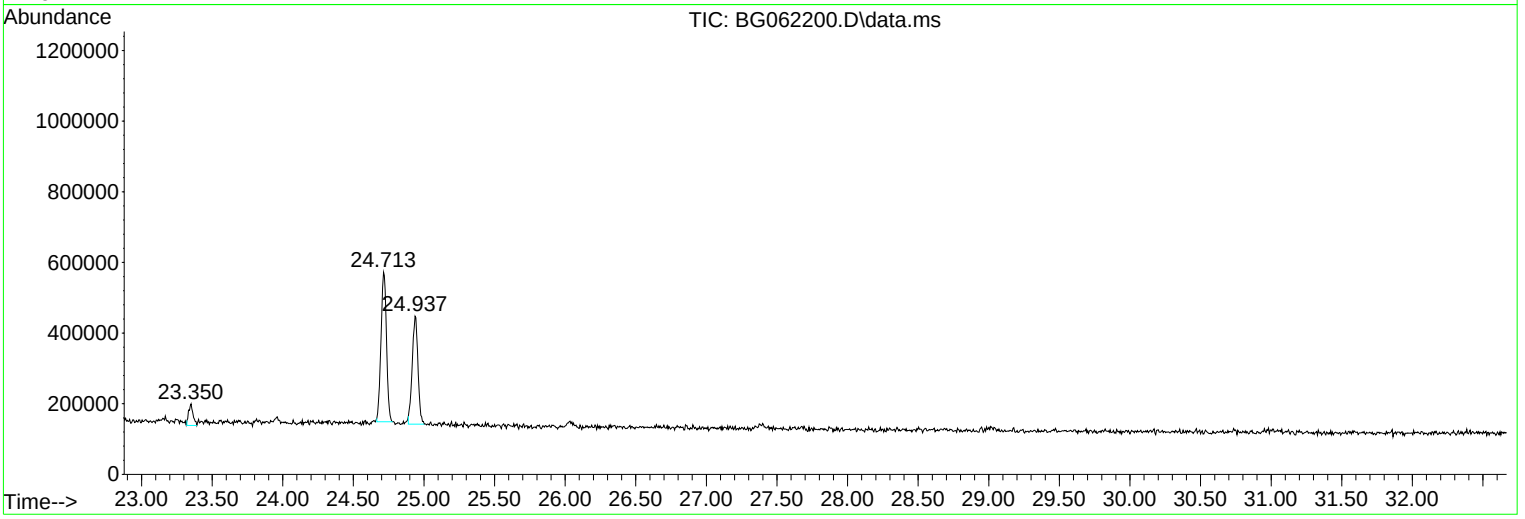
Instrument :
BNA_G
ClientSampleId :
YDZF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

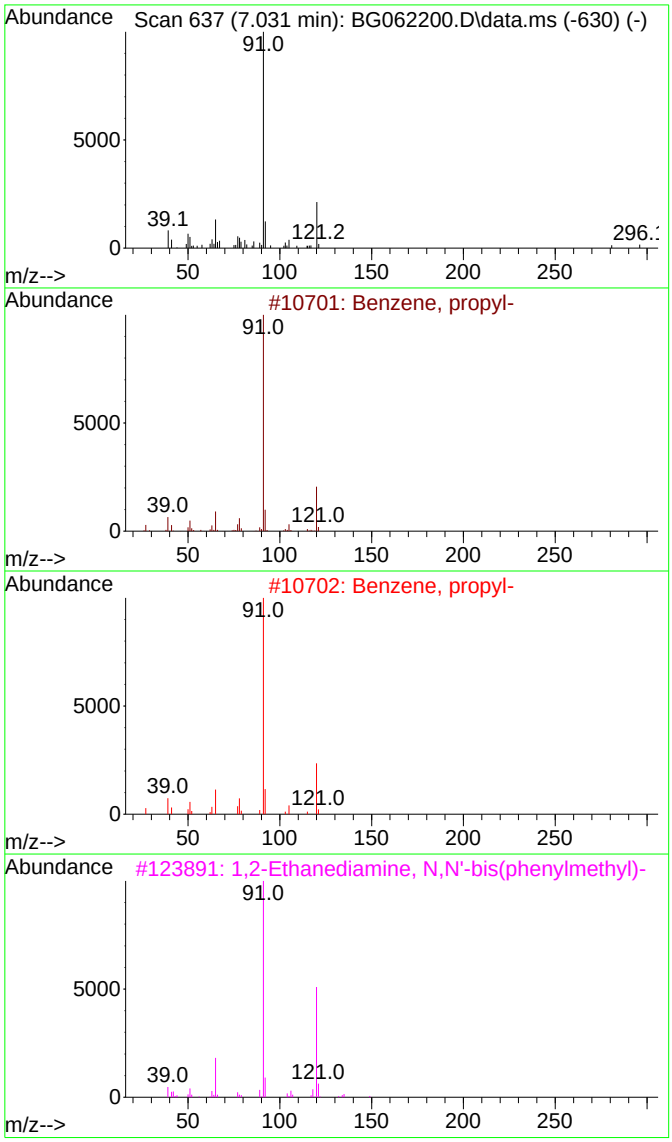
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

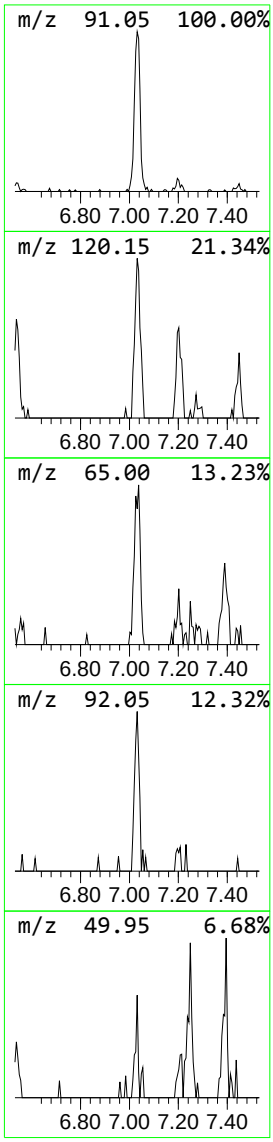
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.031	4.04 ng/ul	66550	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	81
2	Benzene, propyl-	120	C9H12	000103-65-1	74
3	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64
4	1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	64
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

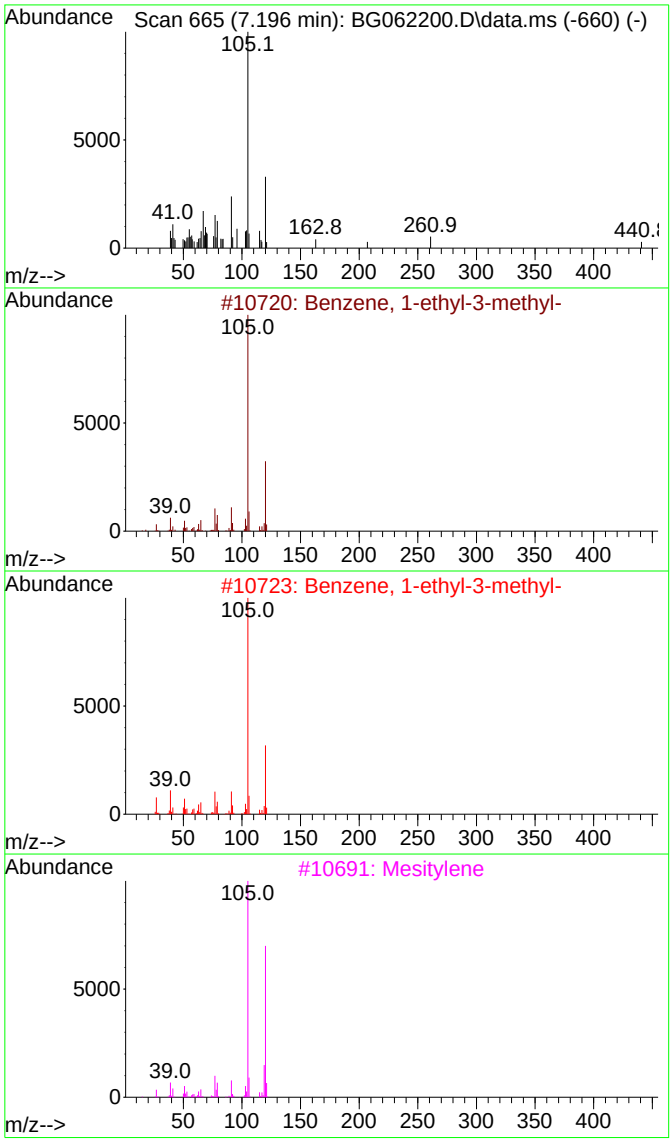
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

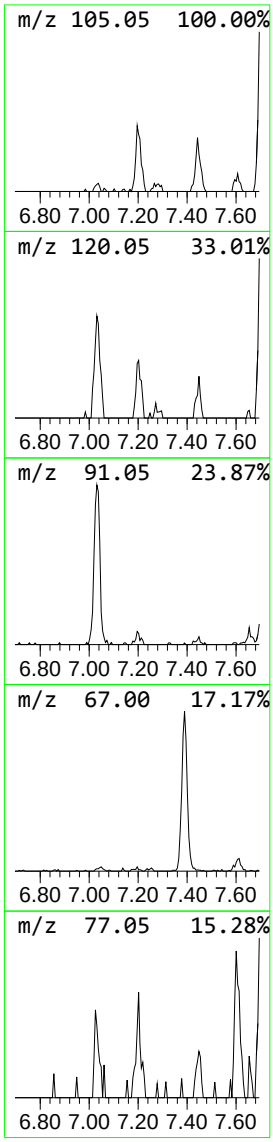
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.196	2.26 ng/ul	37162	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	58
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	58
3	Mesitylene	120	C9H12	000108-67-8	52
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	52
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

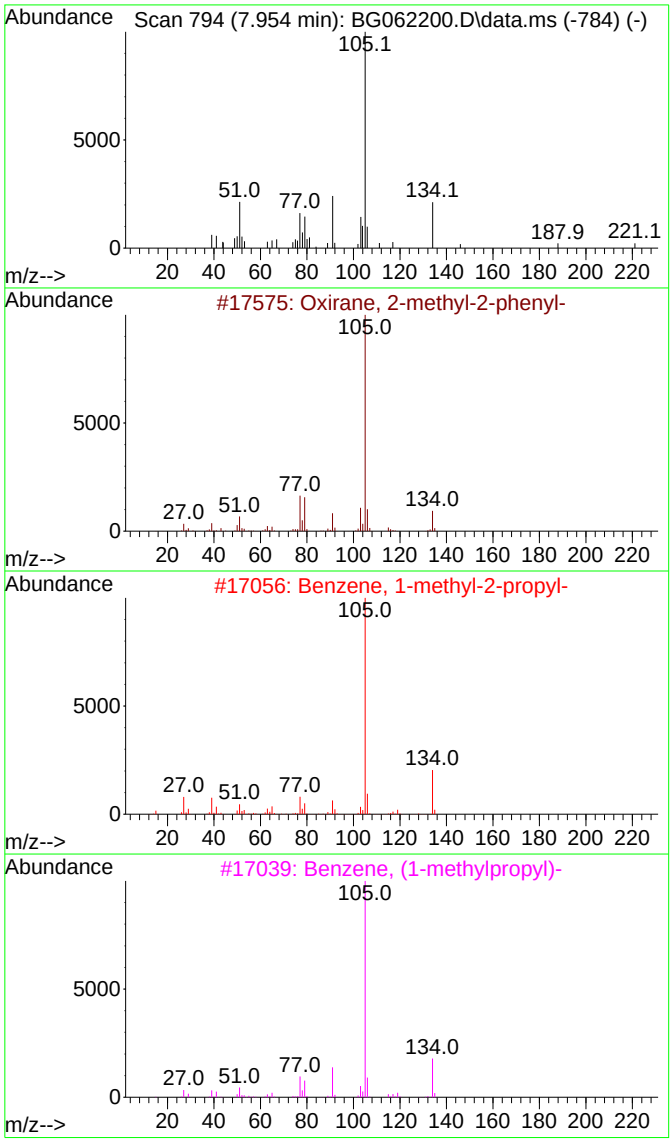
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

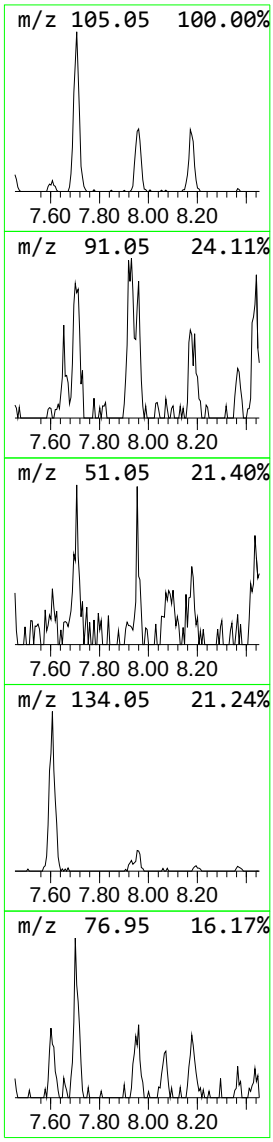
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Oxirane, 2-methyl-2-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.954	3.75 ng/ul	61658	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	74
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	64
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	64
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	58
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

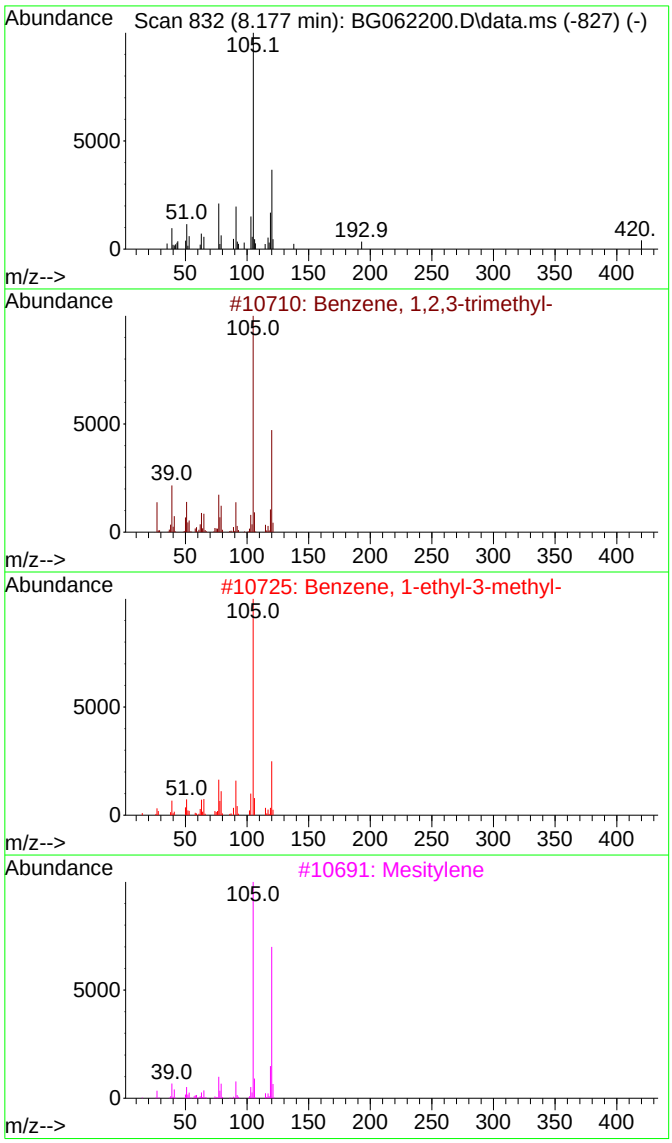
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

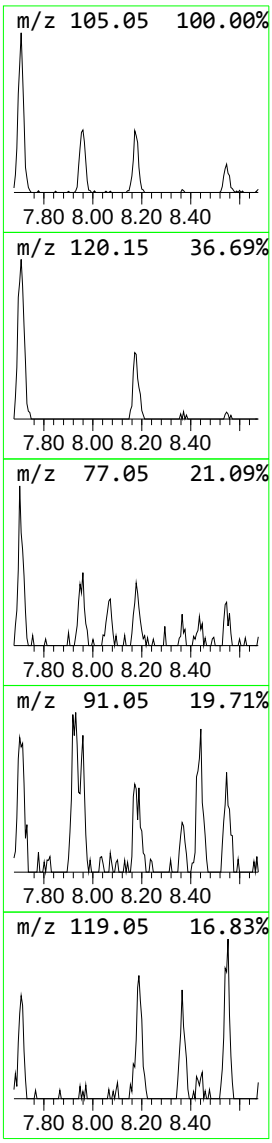
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.177	2.92 ng/ul	48069	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	72
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	72
3	Mesitylene	120	C9H12	000108-67-8	64
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64
5	2,4-Nonadiyne	120	C9H12	063621-15-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

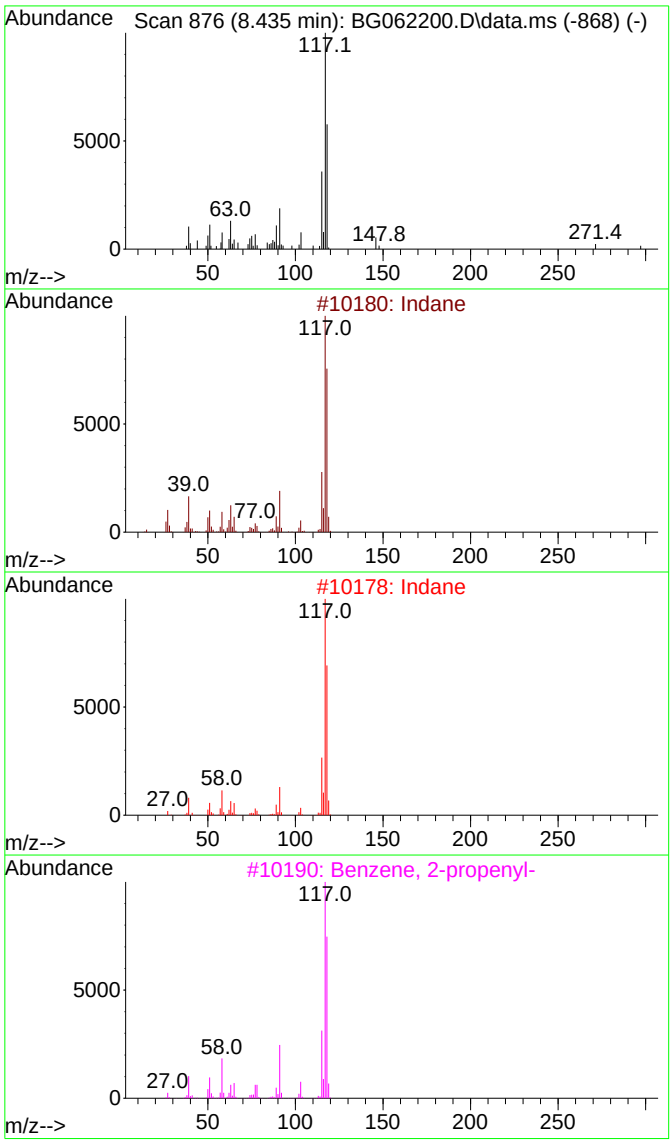
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

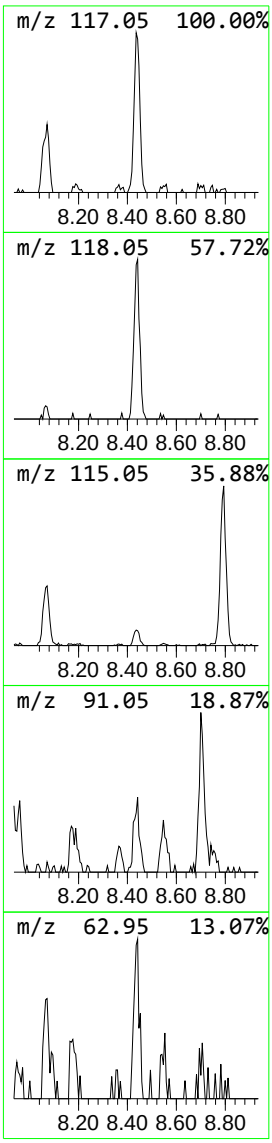
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.435	4.82 ng/ul	79310	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	90
2	Indane	118	C9H10	000496-11-7	87
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	81
5	Deltacyclene	118	C9H10	007785-10-6	81





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

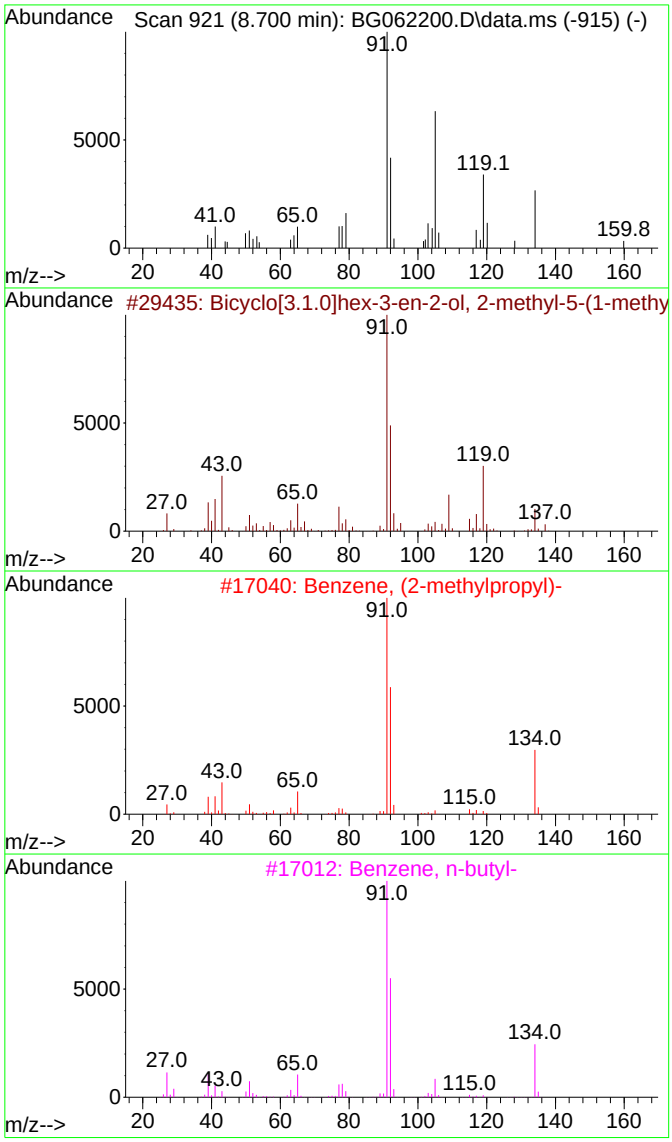
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

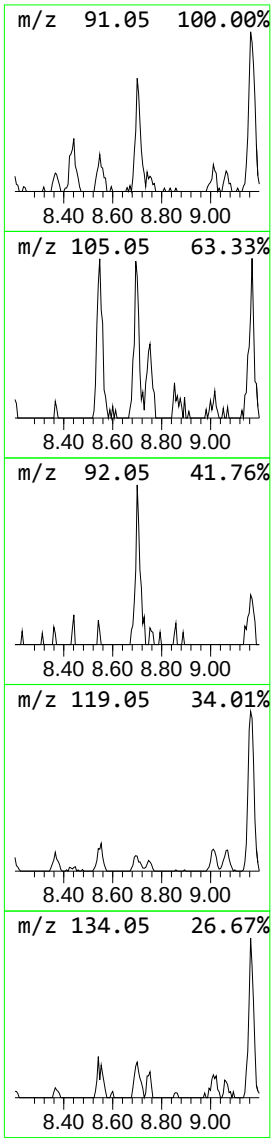
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Bicyclo[3.1.0]hex-3-en-2-ol... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.700	2.64 ng/ul	43416	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.0]hex-3-en-2-ol, 2-m...	152	C10H16O	097631-68-0	50
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	46
3	Benzene, n-butyl-	134	C10H14	000104-51-8	46
4	Benzene, n-butyl-	134	C10H14	000104-51-8	46
5	Benzene, n-butyl-	134	C10H14	000104-51-8	46





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

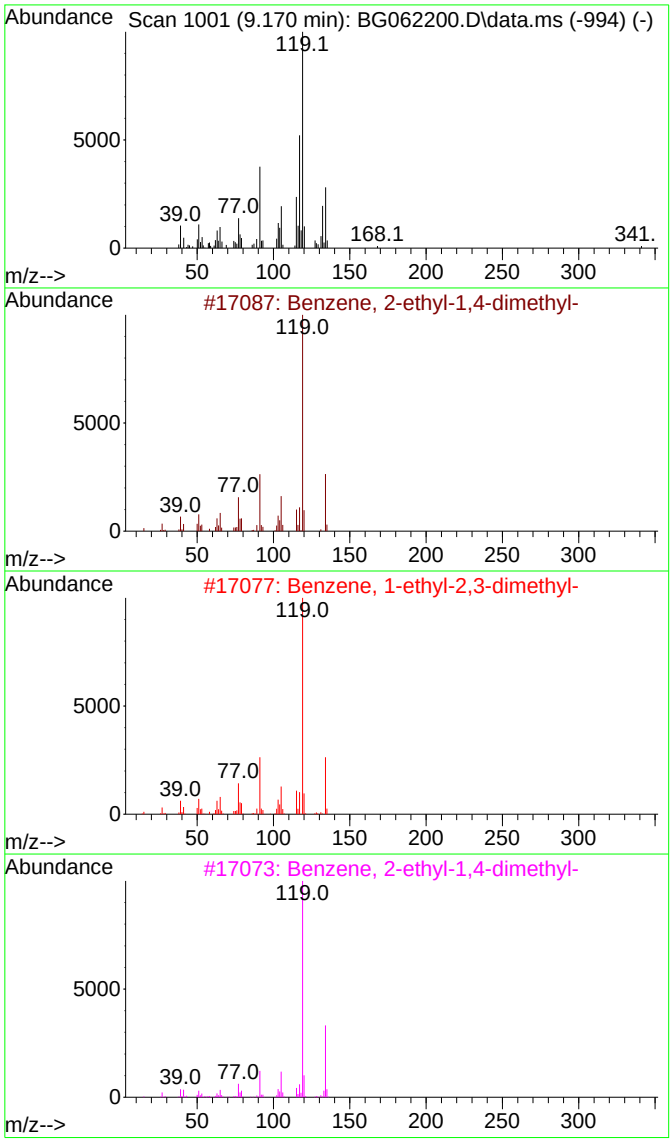
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

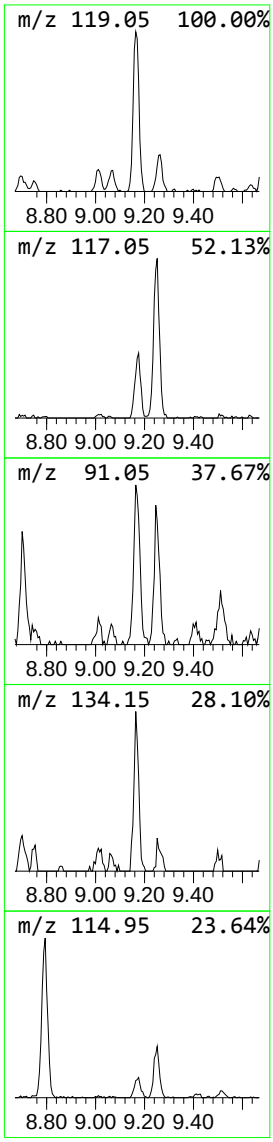
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.170	9.59 ng/ul	157731	1,4-Dichlorobenzene-d4	8.065

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	64
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	58





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Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

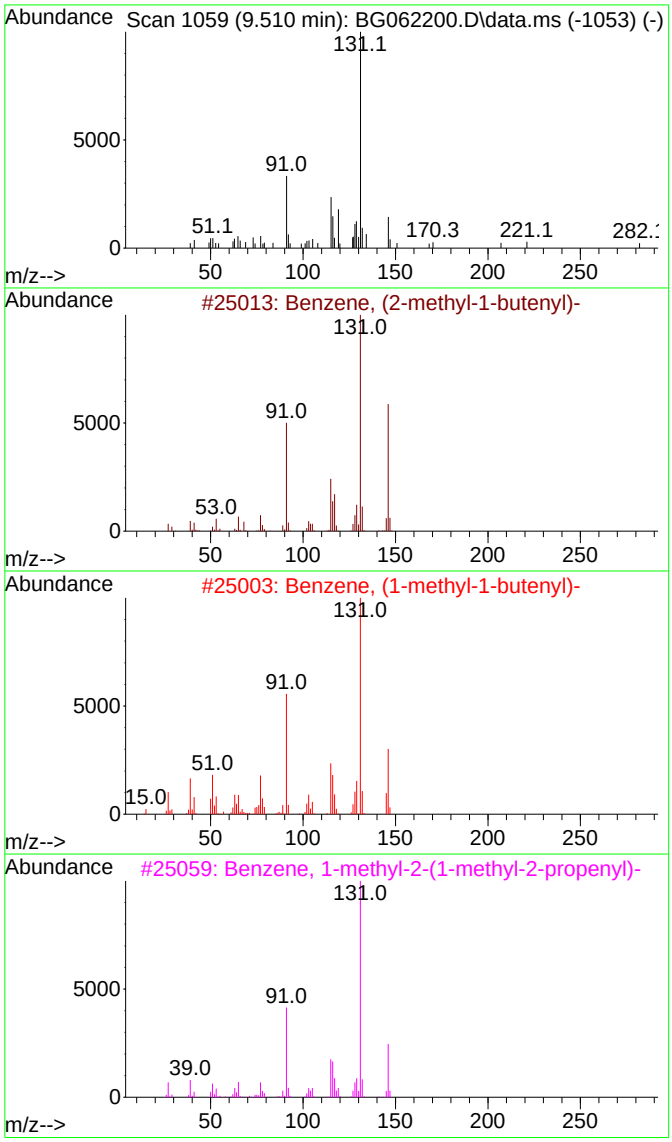
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

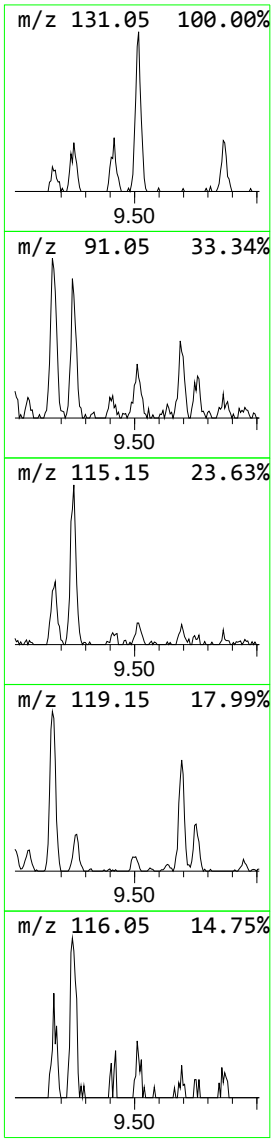
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, (2-methyl-1-butenyl)- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	2.17 ng/ul	52113	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	87
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87
3	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	80
4	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	80
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	80





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

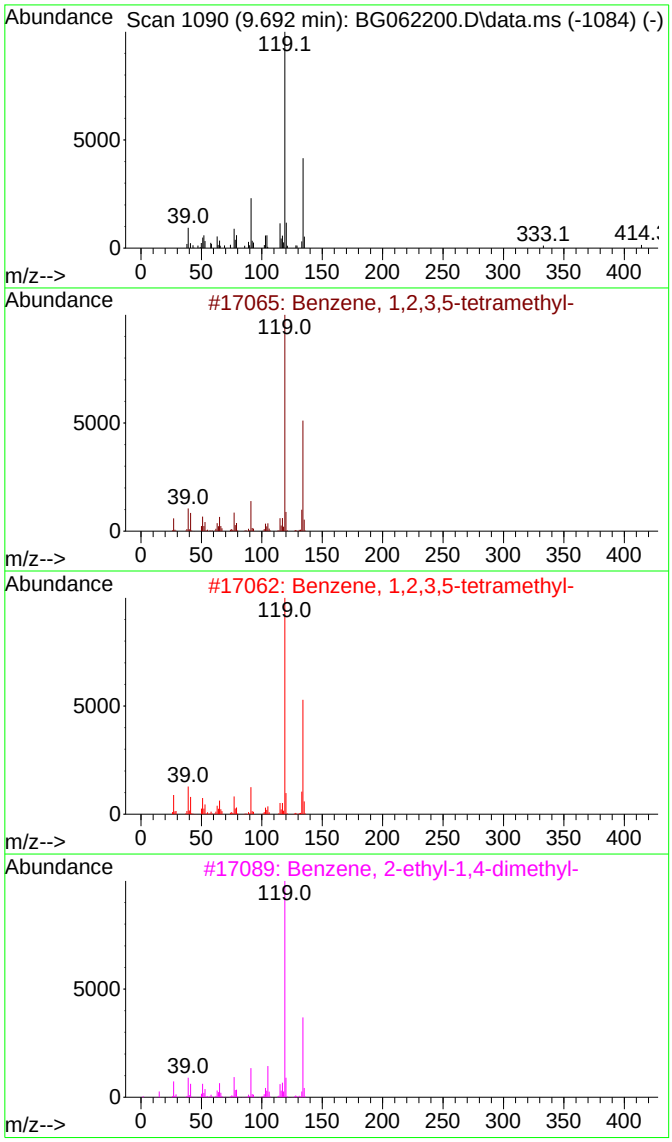
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

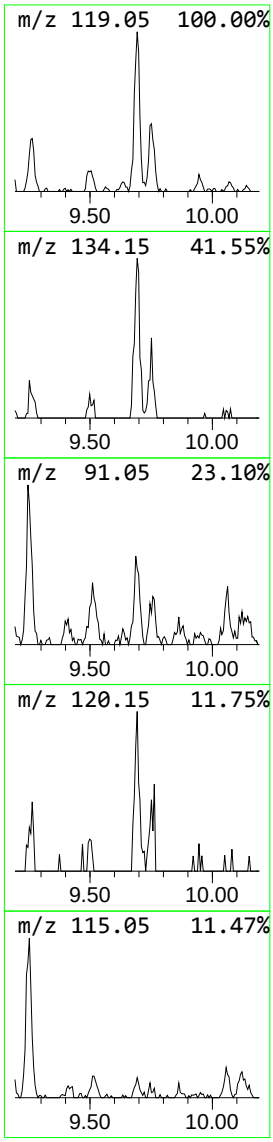
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.692	2.87 ng/ul	68914	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
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Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

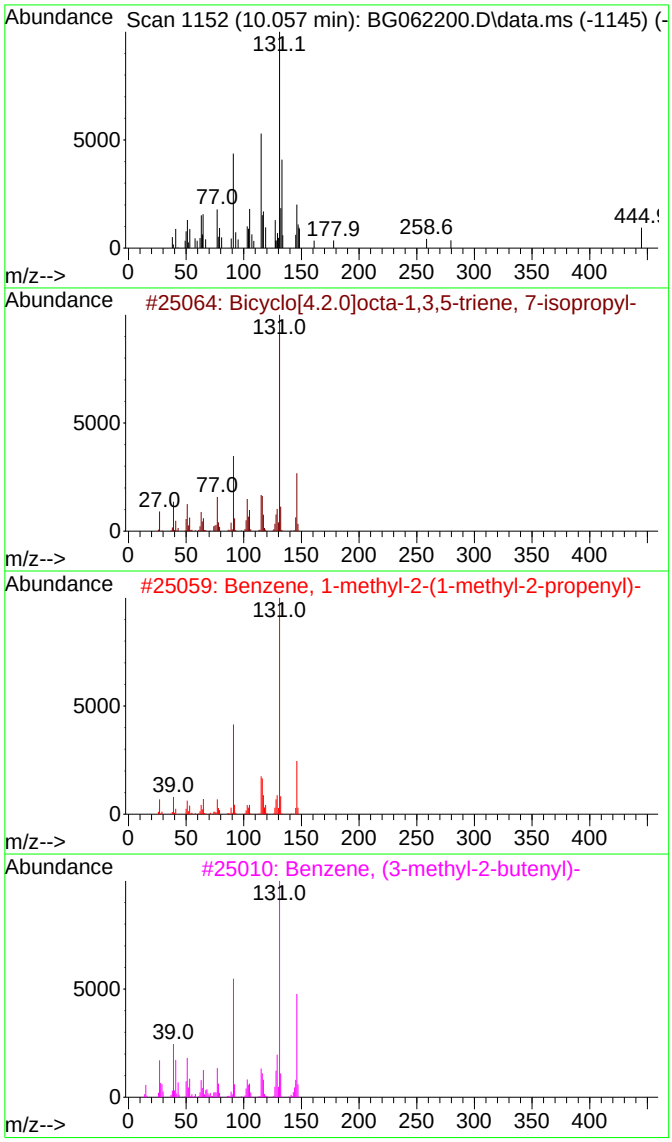
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

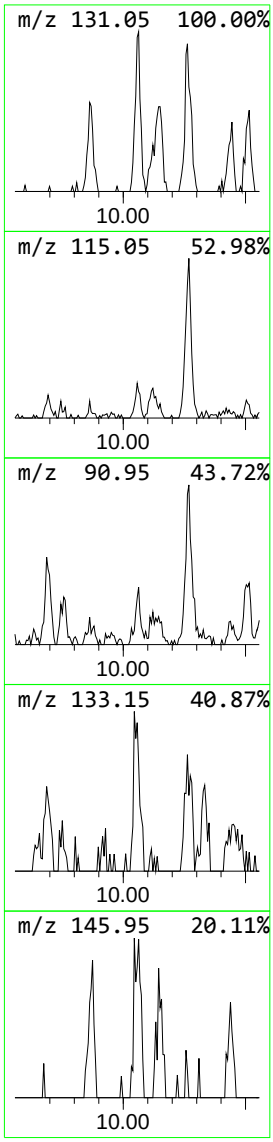
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 unknown-01 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.057	2.02 ng/ul	48475	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	50
2	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	43
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	43
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	38
5	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	38





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

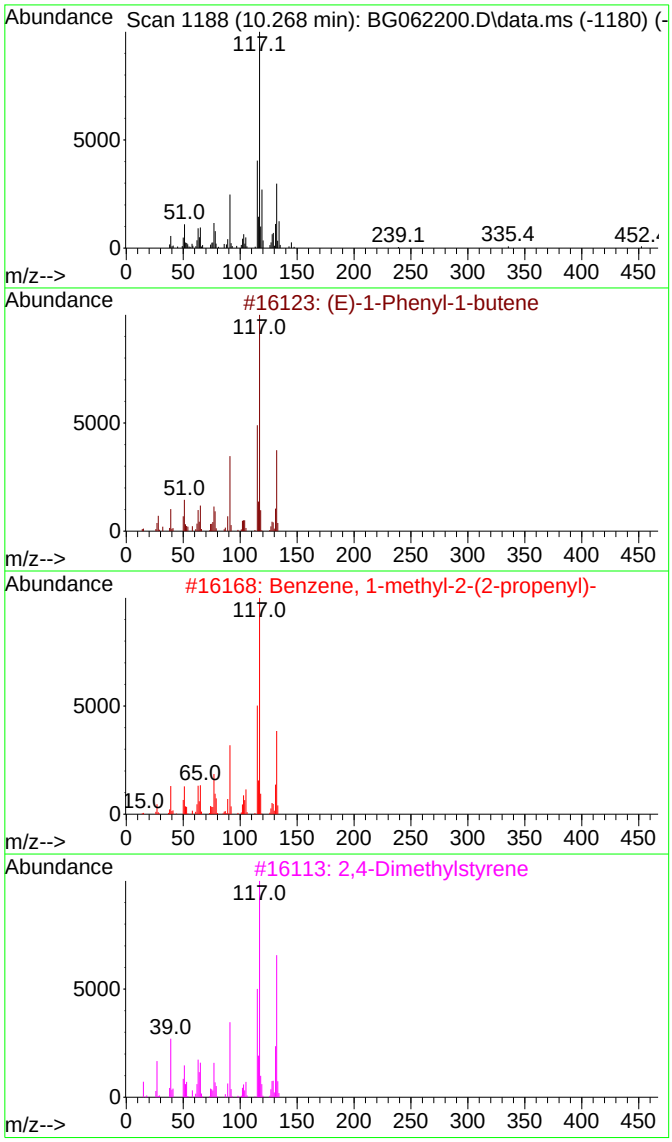
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

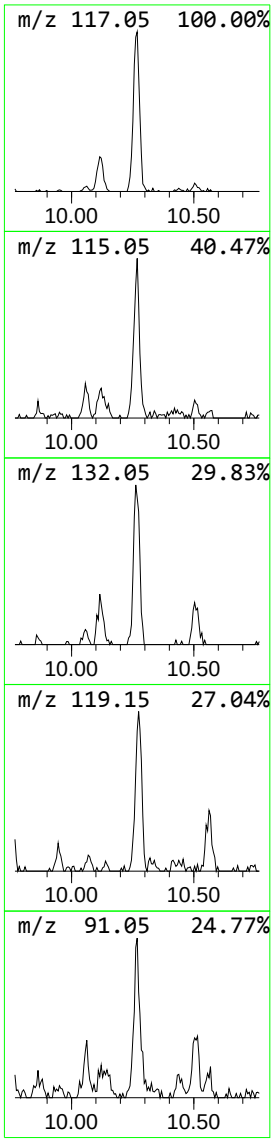
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 (E)-1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.268	8.94 ng/ul	214628	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
3	2,4-Dimethylstyrene	132	C10H12	002234-20-0	90
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

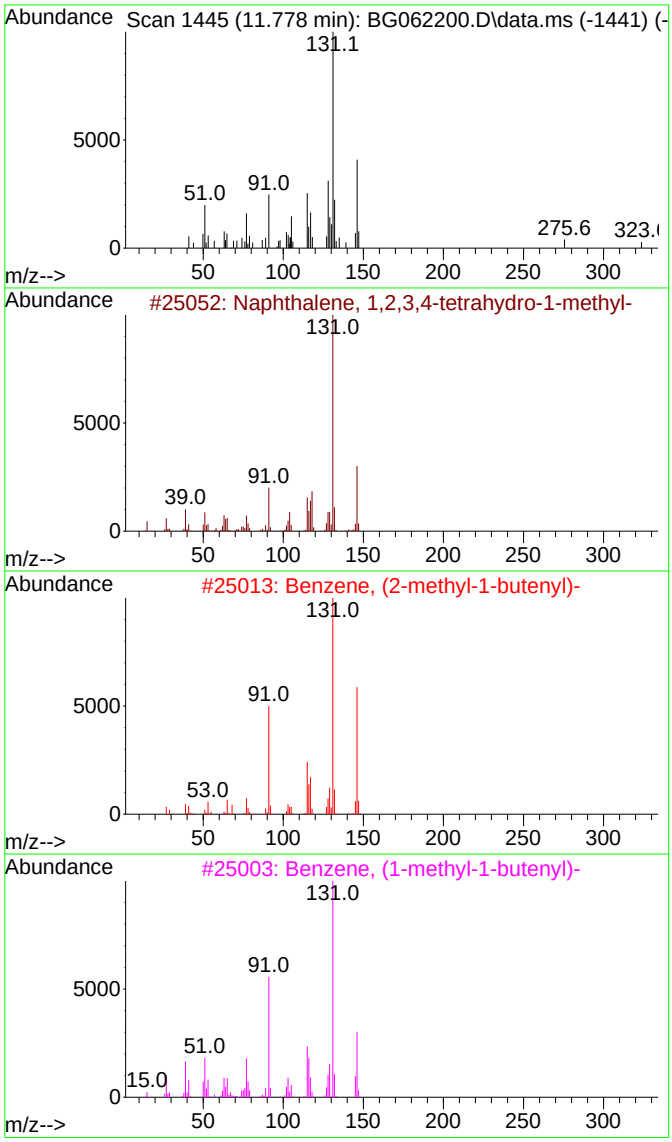
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

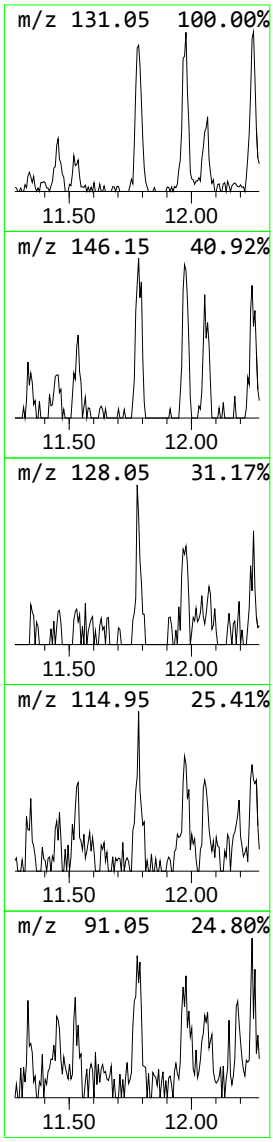
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.778	2.21 ng/ul	53078	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	76
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	74
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	74
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	72





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Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

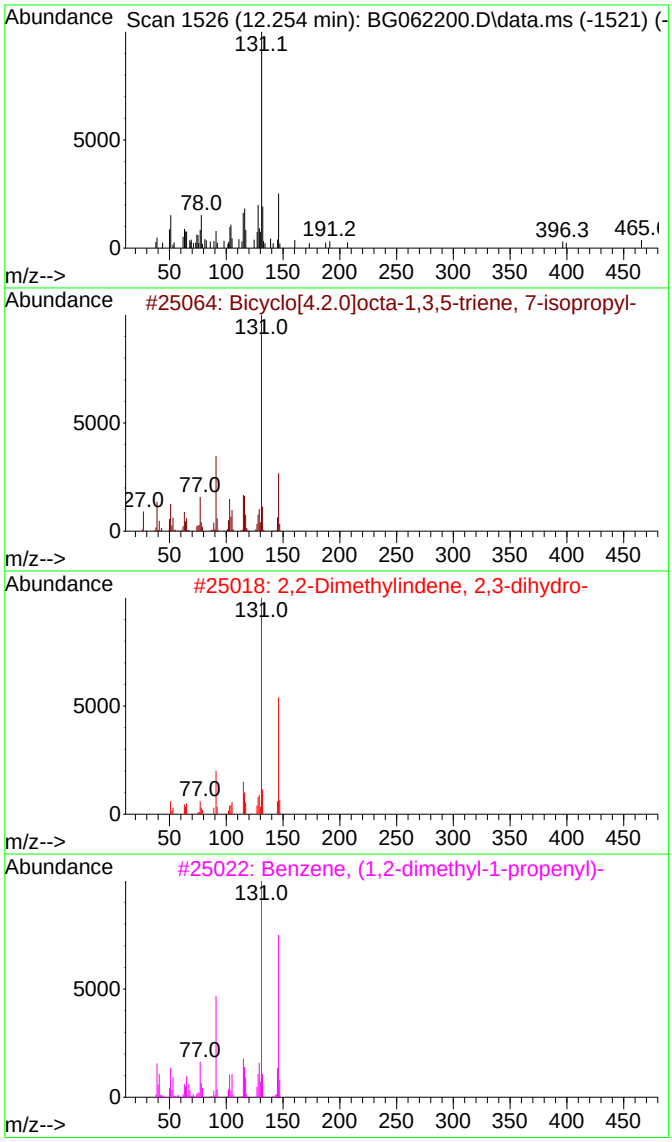
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

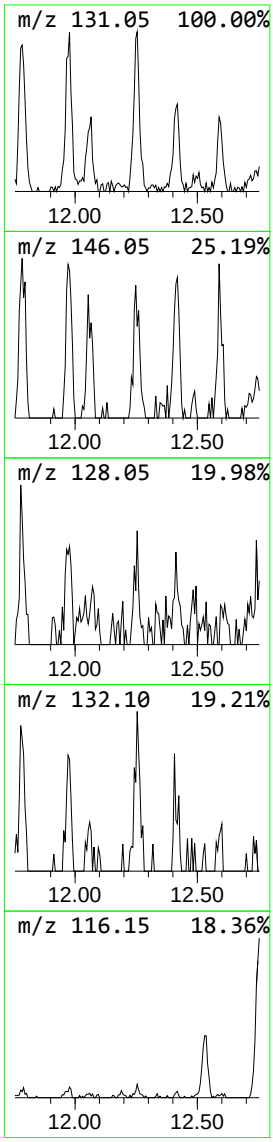
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.254	2.03 ng/ul	48750	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	76
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	74
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	74
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
5	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	055255-94-2	59





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Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

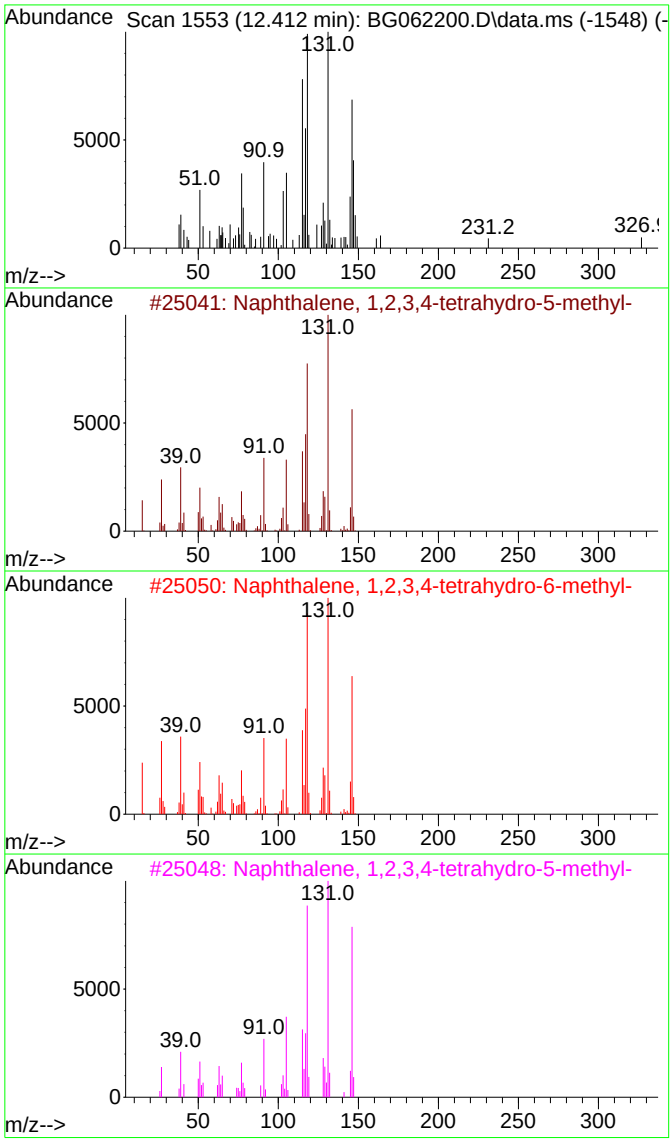
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

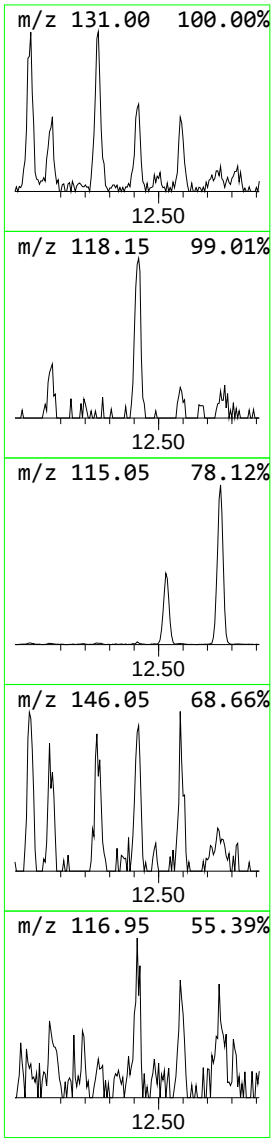
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.412	2.42 ng/ul	57981	Naphthalene-d8	10.885

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	86
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	60
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58





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Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

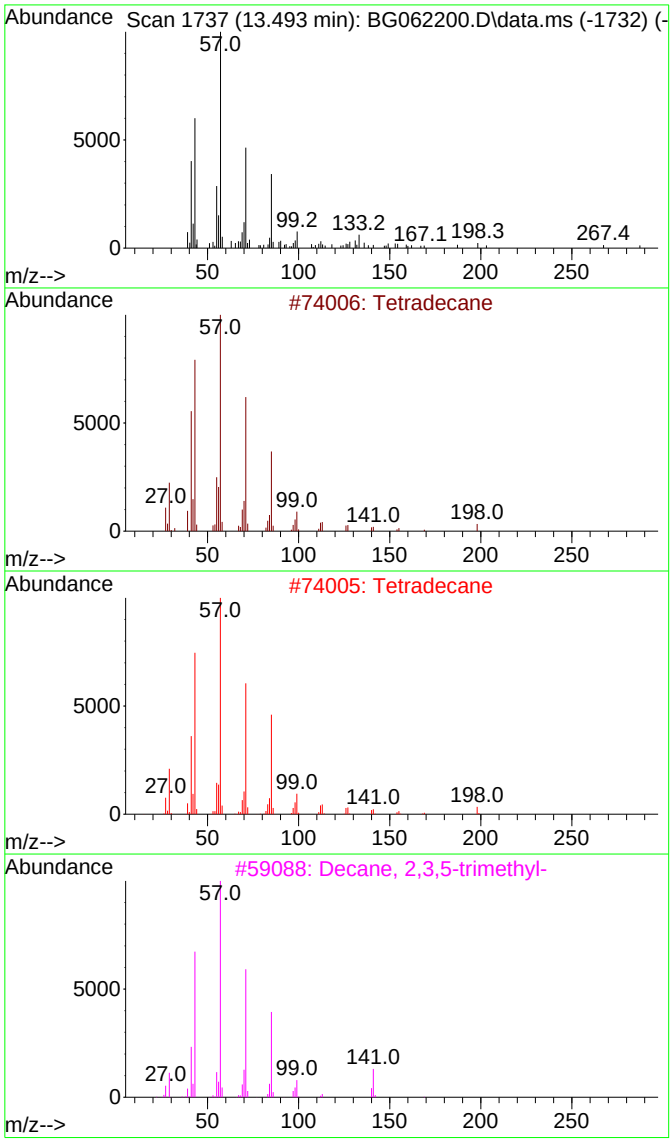
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

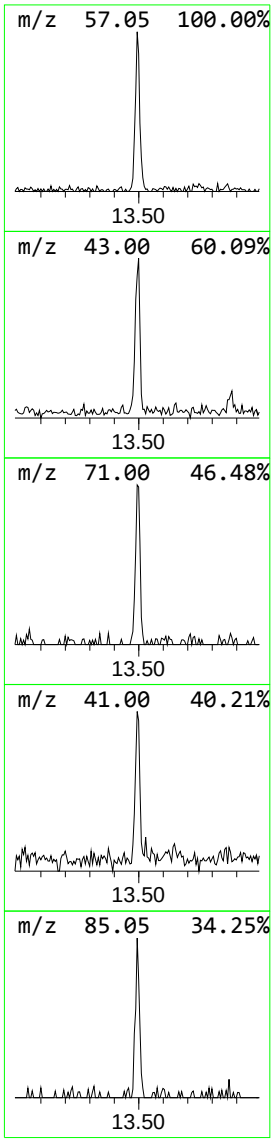
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.493	2.88 ng/ul	101721	Acenaphthene-d10	14.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	89
2	Tetradecane	198	C14H30	000629-59-4	89
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72
4	Hexadecane	226	C16H34	000544-76-3	72
5	Dodecane, 2,5-dimethyl-	198	C14H30	056292-65-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

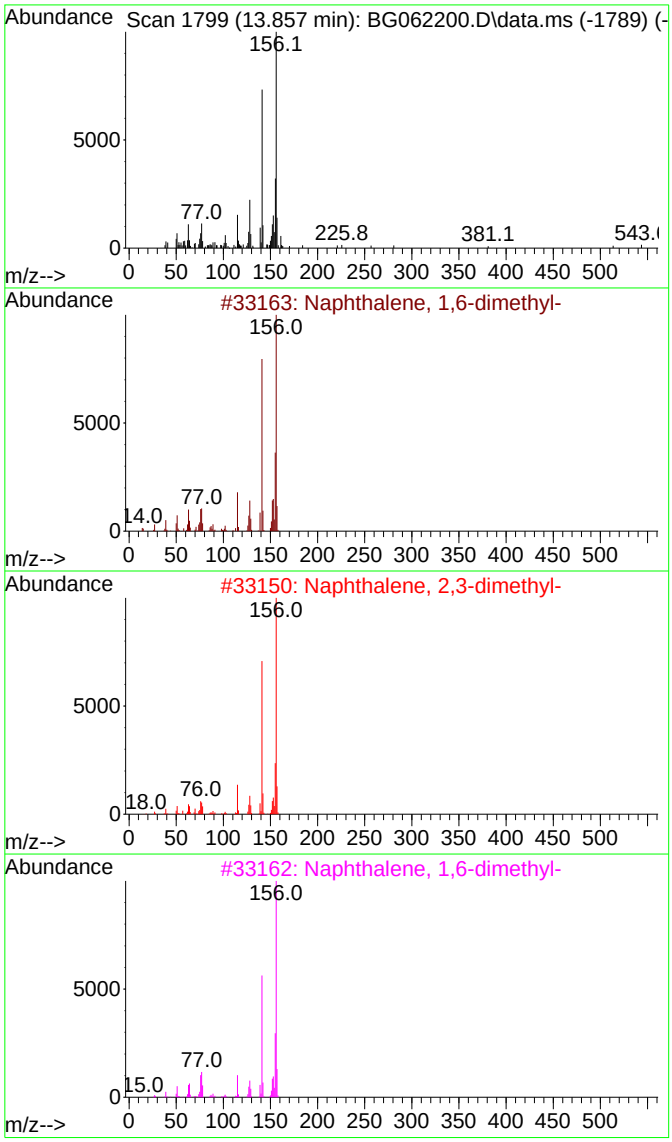
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

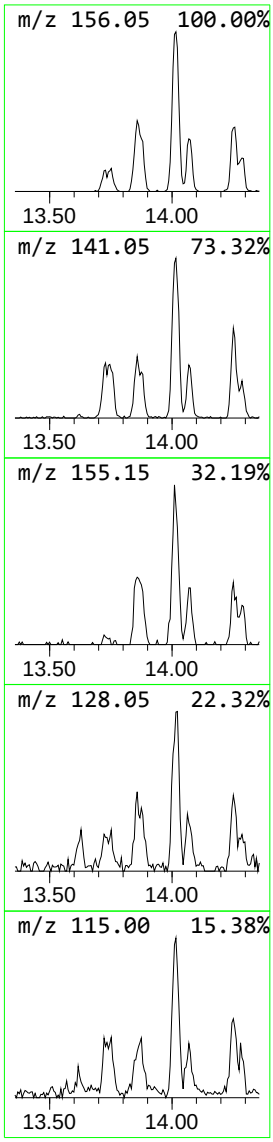
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Naphthalene, 1,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	4.14 ng/ul	146143	Acenaphthene-d10	14.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

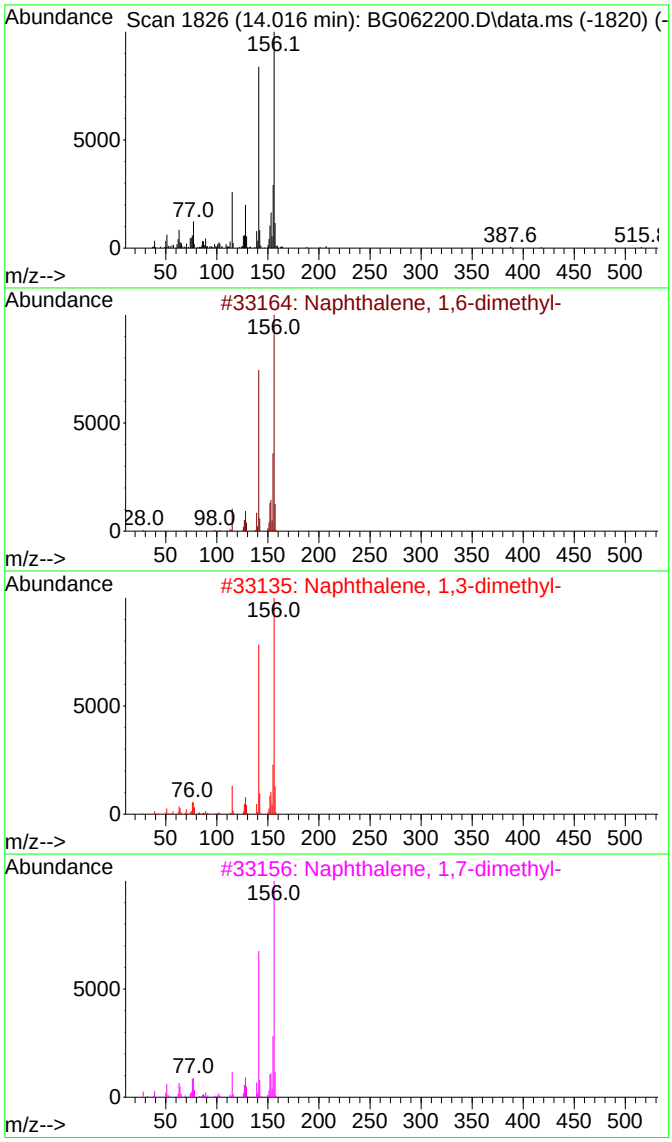
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

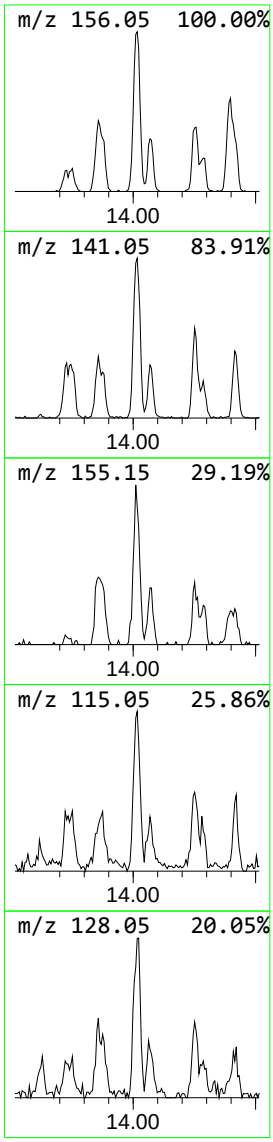
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Naphthalene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.016	9.48 ng/ul	334295	Acenaphthene-d10	14.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	94
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

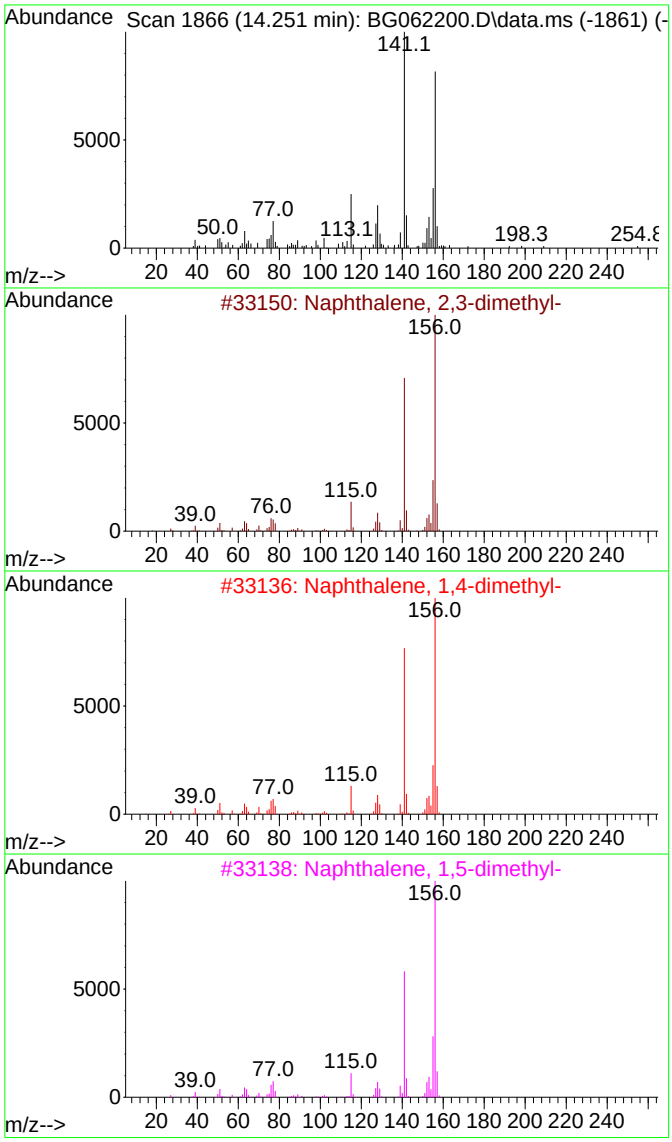
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

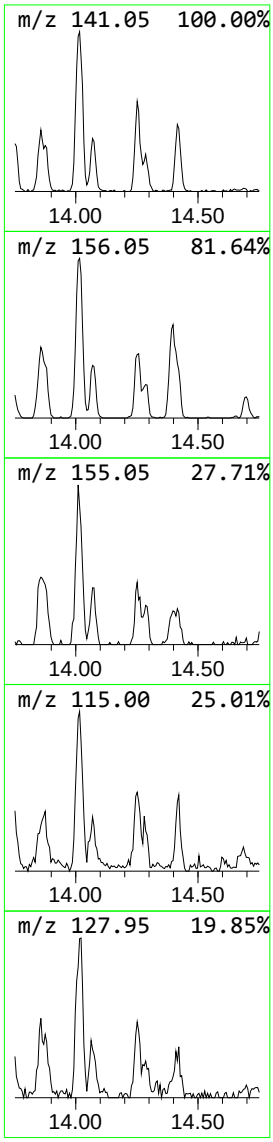
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.251	3.83 ng/ul	135107	Acenaphthene-d10	14.698

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
3	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

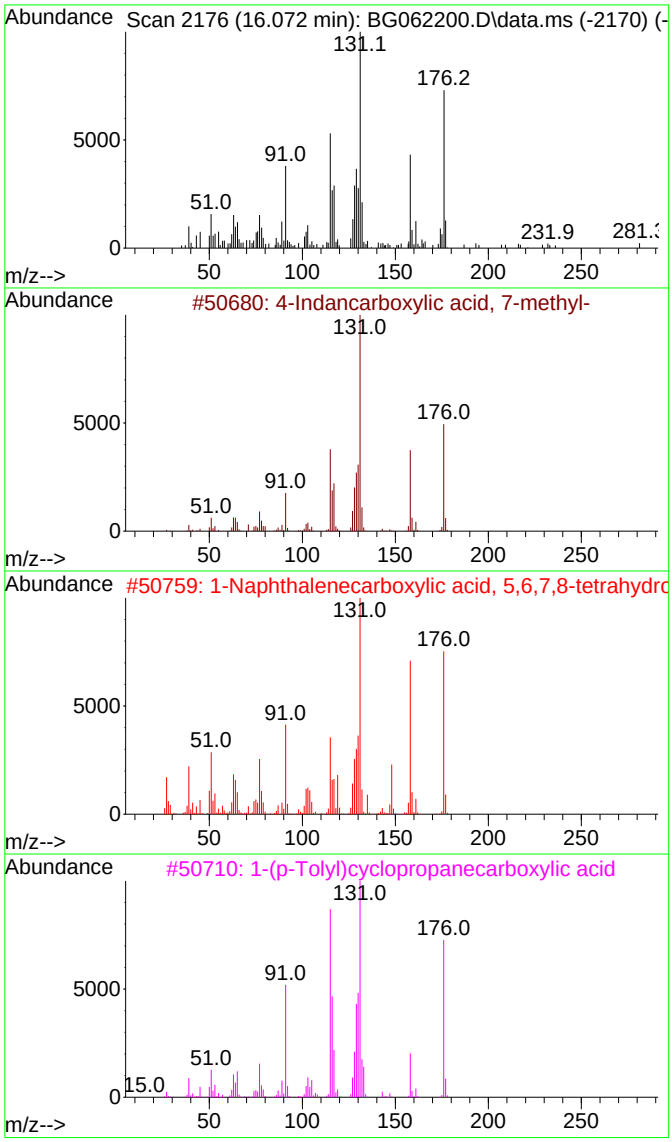
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

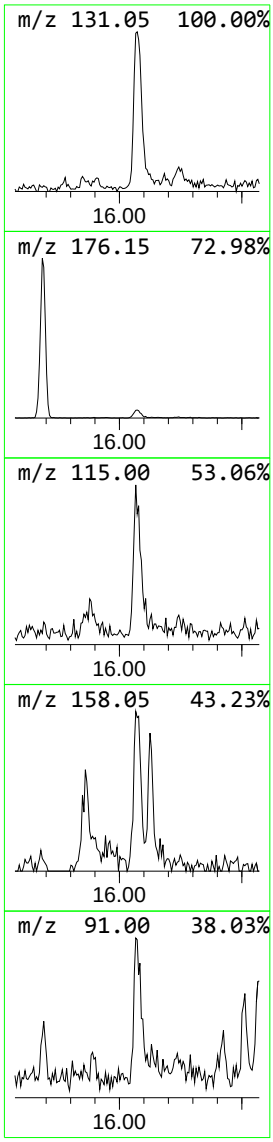
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 4-Indancarboxylic acid, 7-m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.072	5.32 ng/ul	215648	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Indancarboxylic acid, 7-methyl-	176	C11H12O2	071042-74-5	96
2	1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	90
3	1-(p-Tolyl)cyclopropanecarboxyli...	176	C11H12O2	083846-66-6	64
4	2-(1,2,3,4-Tetrahydronaphthalen-...	176	C12H16O	068480-12-6	52
5	2,3-Dihydro-1H-inden-5-ylacetic ...	176	C11H12O2	005453-98-5	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

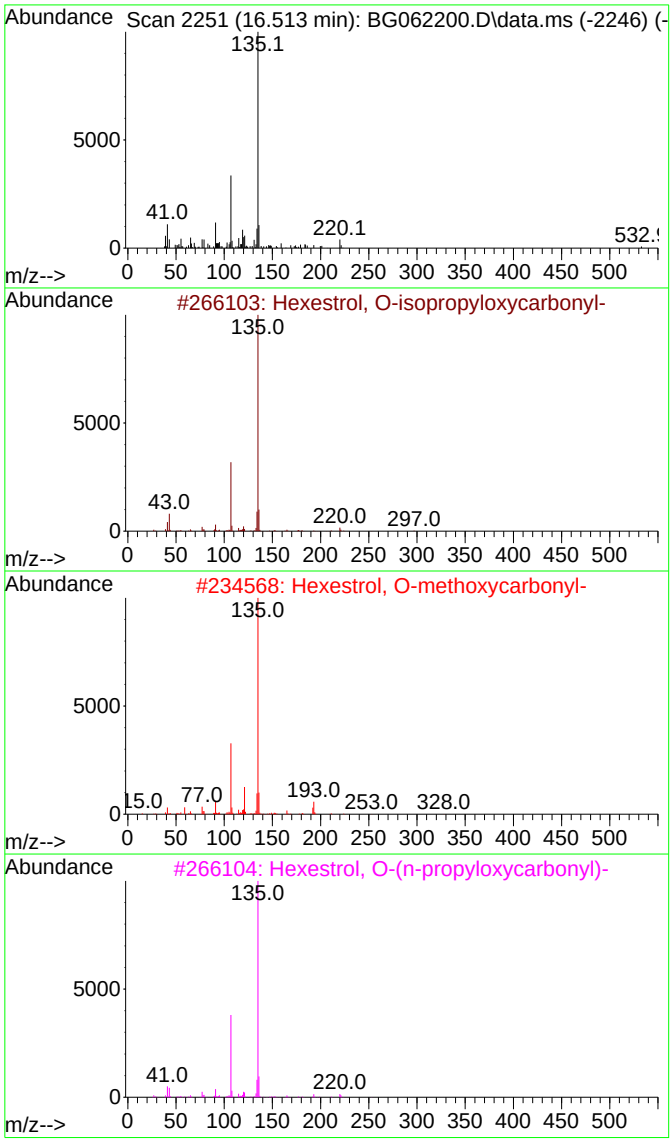
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

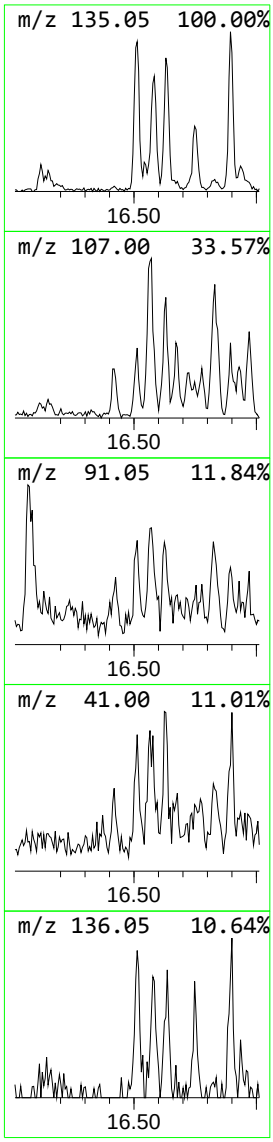
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Hexestrol, O-isopropyloxyca... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.513	2.50 ng/ul	101055	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexestrol, O-isopropyloxycarbonyl-	356	C22H28O4	1000487-68-4	72
2	Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	72
3	Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	72
4	Hexestrol, O-acetyl-	312	C20H24O3	1000374-87-8	72
5	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

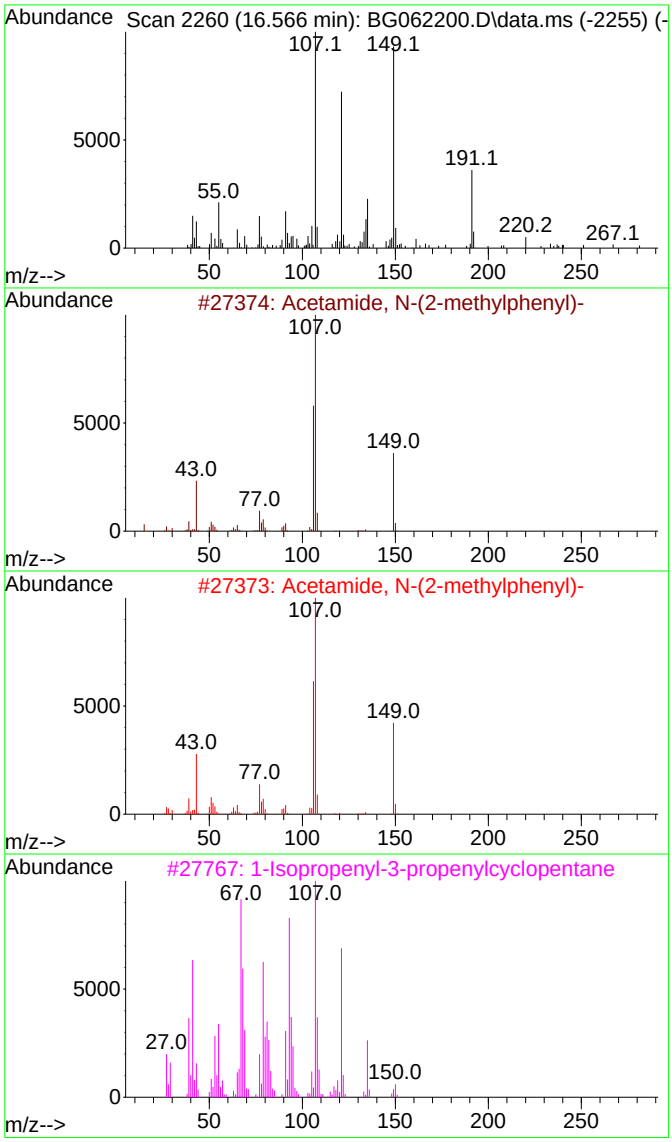
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

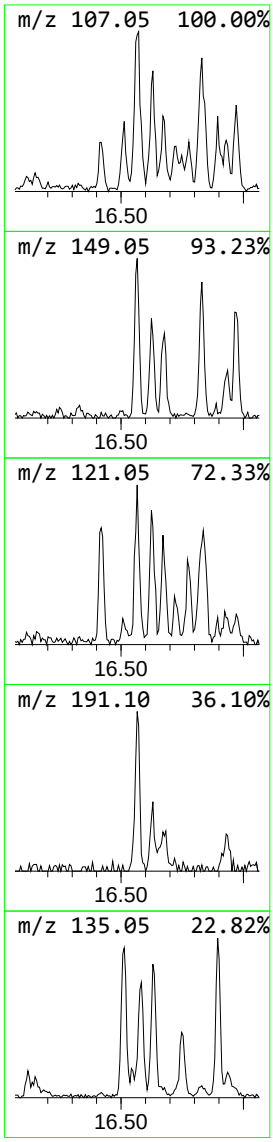
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Acetamide, N-(2-methylphenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.566	4.99 ng/ul	202069	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	53
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
3	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	43
4	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
5	Benzene, 1-ethoxy-4-ethyl-	150	C10H14O	001585-06-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

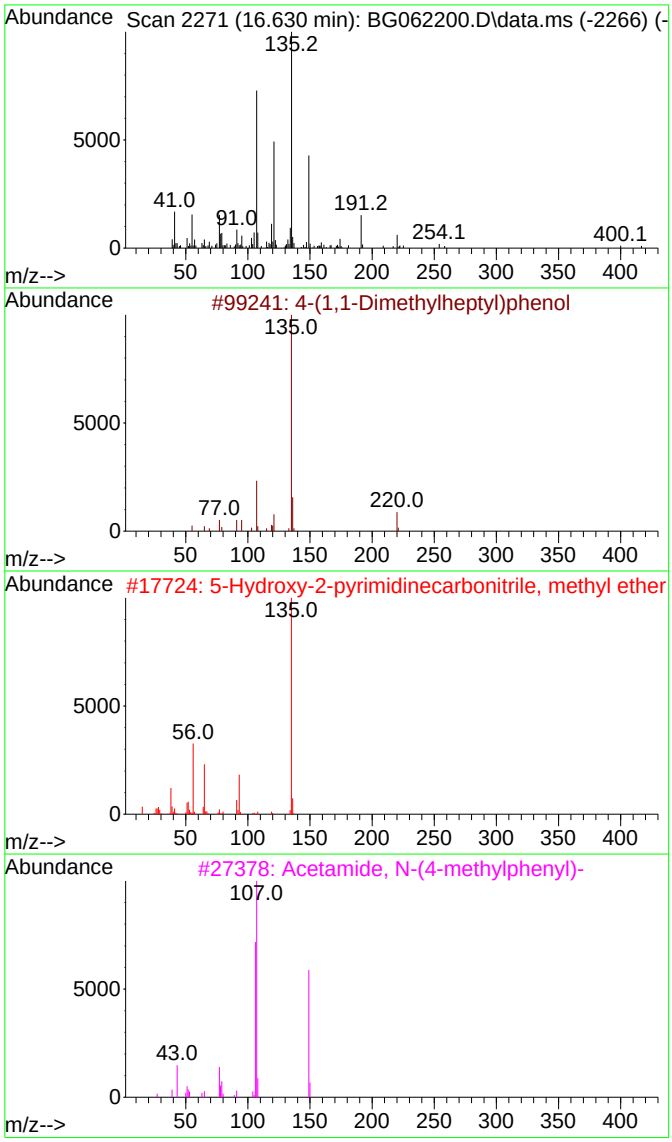
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

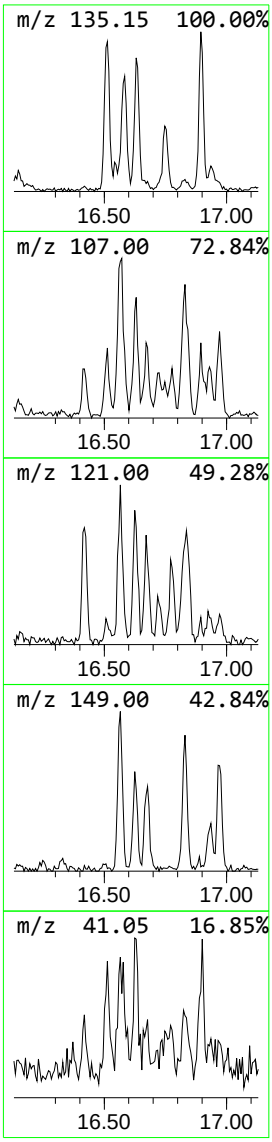
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 4-(1,1-Dimethylheptyl)phenol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.630	3.54 ng/ul	143537	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	50
2	5-Hydroxy-2-pyrimidinecarbonitri...	135	C6H5N3O	087362-32-1	27
3	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	25
4	Acetamide, N-methyl-N-phenyl-	149	C9H11NO	000579-10-2	25
5	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

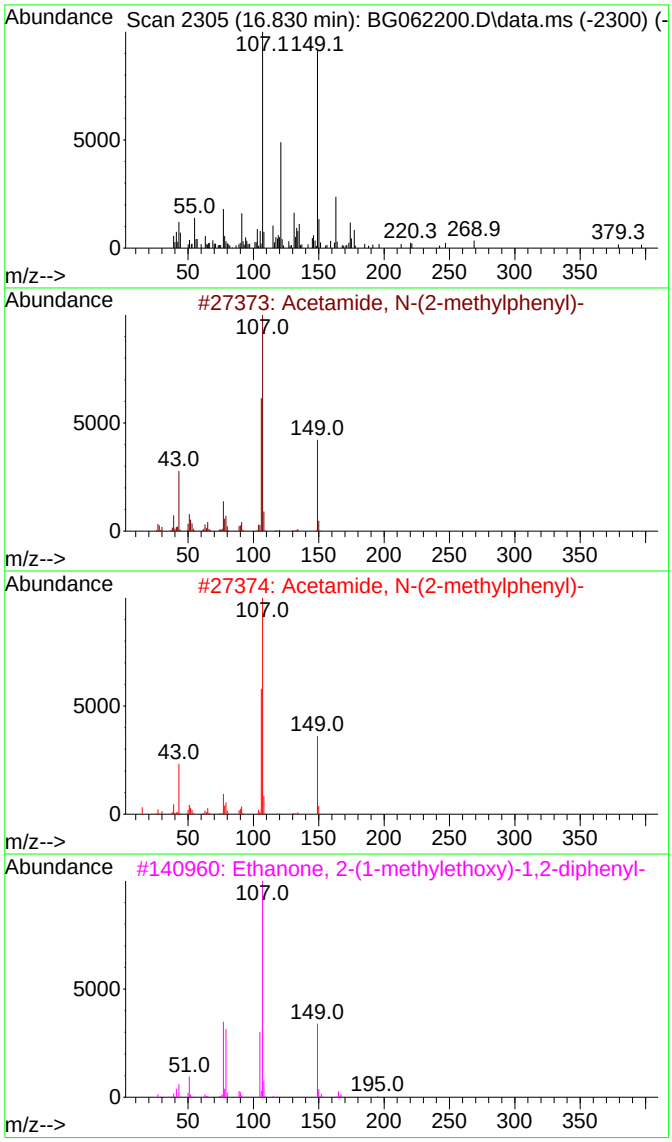
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

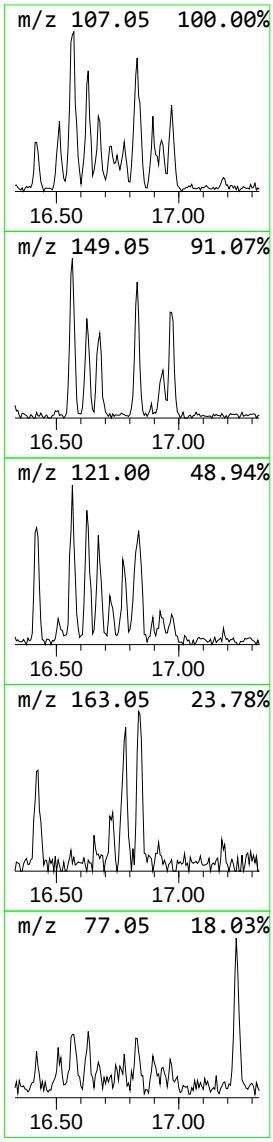
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Ethanone, 2-(1-methylethoxy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.830	3.79 ng/ul	153551	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	58
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	53
3	Ethanone, 2-(1-methylethoxy)-1,2...	254	C17H18O2	006652-28-4	53
4	(S)-5-((3-Acetylphenoxy)methyl)-...	379	C21H21N3O4	1623481-80-0	38
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

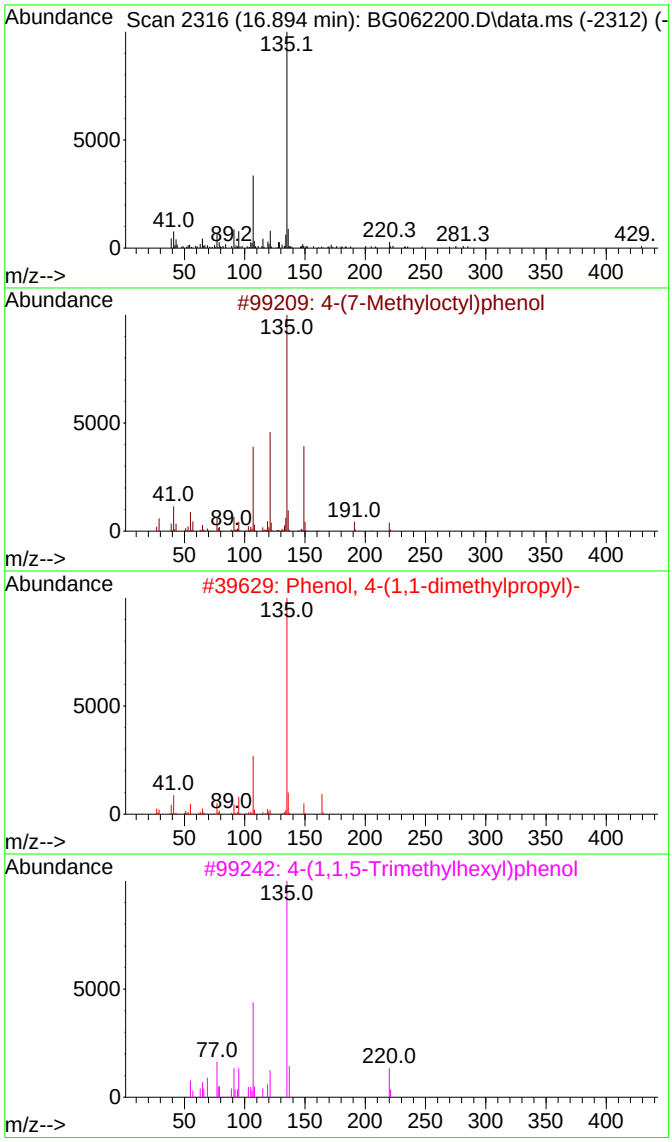
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

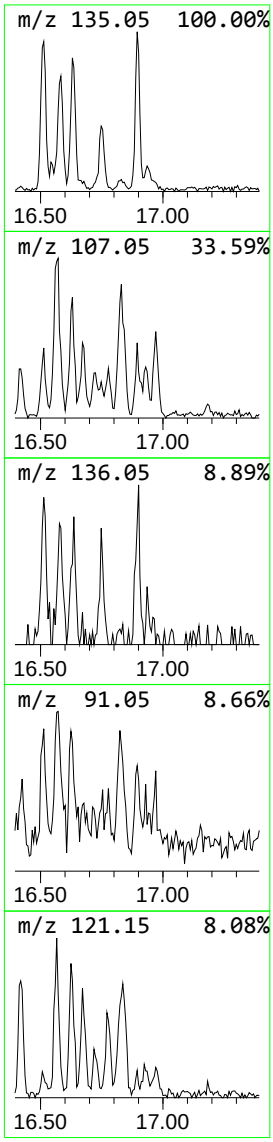
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 4-(7-Methyloctyl)phenol Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.895	2.04 ng/ul	82437	Phenanthrene-d10	17.441

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	87
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	80
3	4-(1,1,5-Trimethylhexyl)phenol	220	C15H24O	1000384-80-1	74
4	Hexestrol	270	C18H22O2	000084-16-2	72
5	Imidazo[1,2-c]pyrimidin-5-ol	135	C6H5N3O	055662-66-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

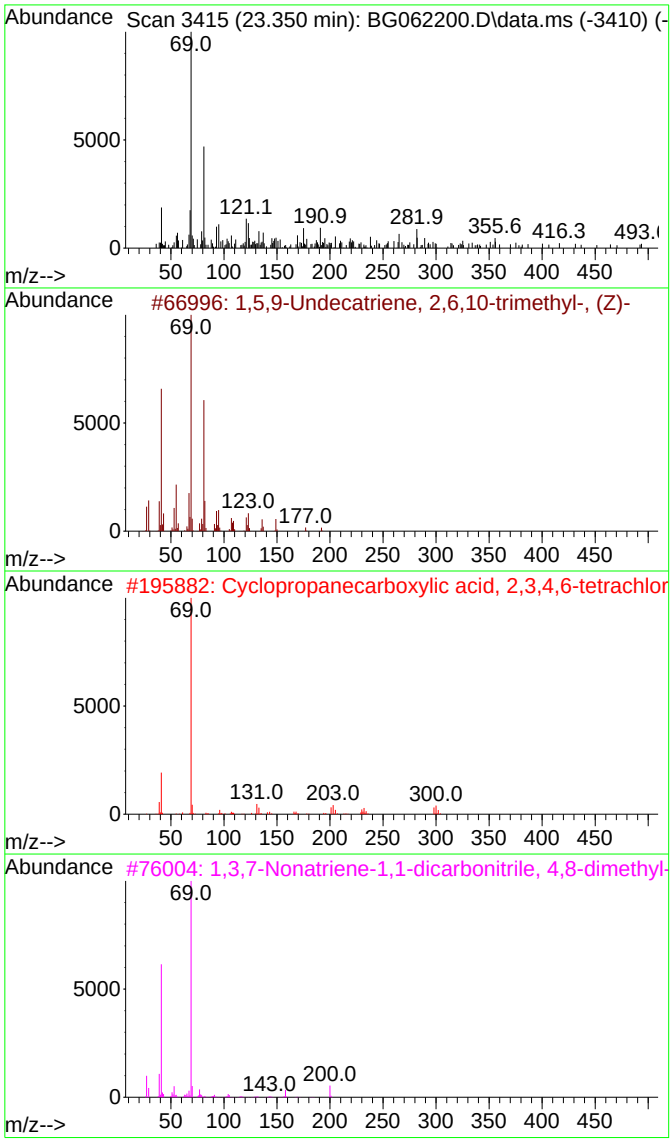
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

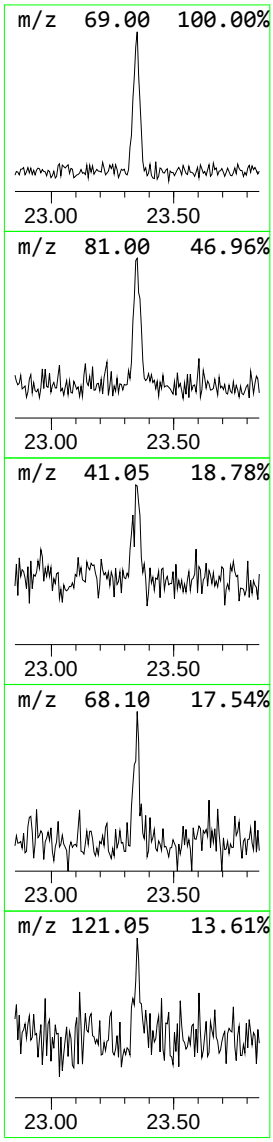
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 1,5,9-Undecatriene, 2,6,10-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.351	3.01 ng/ul	129915	Perylene-d12	24.936

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	58
2	Cyclopropanecarboxylic acid, 2,3...	298	C10H6Cl4O2	1000354-66-5	53
3	1,3,7-Nonatriene-1,1-dicarbonitr...	200	C13H16N2	1000162-13-4	53
4	3,6-Dimethyl-1-heptyn-3-ol	140	C9H16O	019549-98-5	53
5	4-Hexenenitrile, 2,5-dimethyl-2-...	191	C13H21N	069340-56-3	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062200.D
Acq On : 24 Jul 2024 00:44
Operator : MA/JU
Sample : P3244-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YDZF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, propyl-	7.031	4.0	ng/ul	66550	1	8.065	329102	20.0
Benzene, 1-ethy...	7.196	2.3	ng/ul	37162	1	8.065	329102	20.0
Oxirane, 2-meth...	7.954	3.8	ng/ul	61658	1	8.065	329102	20.0
Benzene, 1,2,3-...	8.177	2.9	ng/ul	48069	1	8.065	329102	20.0
Indane	8.435	4.8	ng/ul	79310	1	8.065	329102	20.0
Bicyclo[3.1.0]h...	8.700	2.6	ng/ul	43416	1	8.065	329102	20.0
Benzene, 2-ethy...	9.170	9.6	ng/ul	157731	1	8.065	329102	20.0
Benzene, (2-met...	9.510	2.2	ng/ul	52113	2	10.885	480118	20.0
Benzene, 1,2,3,...	9.692	2.9	ng/ul	68914	2	10.885	480118	20.0
unknown-01	10.057	2.0	ng/ul	48475	2	10.885	480118	20.0
(E)-1-Phenyl-1-...	10.268	8.9	ng/ul	214628	2	10.885	480118	20.0
Naphthalene, 1,...	11.778	2.2	ng/ul	53078	2	10.885	480118	20.0
Bicyclo[4.2.0]o...	12.254	2.0	ng/ul	48750	2	10.885	480118	20.0
Naphthalene, 1,...	12.412	2.4	ng/ul	57981	2	10.885	480118	20.0
(DEL) Alkane: S...	13.493	2.9	ng/ul	101721	3	14.698	705243	20.0
Naphthalene, 1,...	13.857	4.1	ng/ul	146143	3	14.698	705243	20.0
Naphthalene, 1,...	14.016	9.5	ng/ul	334295	3	14.698	705243	20.0
Naphthalene, 2,...	14.251	3.8	ng/ul	135107	3	14.698	705243	20.0
4-Indancarboxyl...	16.072	5.3	ng/ul	215648	4	17.441	810060	20.0
Hexestrol, O-is...	16.513	2.5	ng/ul	101055	4	17.441	810060	20.0
Acetamide, N-(2...	16.566	5.0	ng/ul	202069	4	17.441	810060	20.0
4-(1,1-Dimethyl...	16.630	3.5	ng/ul	143537	4	17.441	810060	20.0
Ethanone, 2-(1-...	16.830	3.8	ng/ul	153551	4	17.441	810060	20.0
4-(7-Methylocty...	16.895	2.0	ng/ul	82437	4	17.441	810060	20.0
1,5,9-Undecatri...	23.351	3.0	ng/ul	129915	6	24.936	863126	20.0