LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

Signal : TIC: BG063064.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.487	63	68	71	rBV4	7564	8861	0.33%	0.058%
2	4.057	162	165	168	rBV3	10971	9075	0.34%	0.060%
3	4.322	205	210	216	rVB4	13189	23272	0.88%	0.153%
4	5.379	383	390	393	rVB4	6521	14014	0.53%	0.092%
5	5.497	408	410	413	rBV2	9059	13403	0.51%	0.088%
6	5.873	469	474	481	rVB5	15723	27533	1.04%	0.181%
7	6.354	551	556	558	rBV4	9507	12701	0.48%	0.084%
8	6.725	613	619	623	rBV3	4844	8062	0.30%	0.053%
9	6.836	632	638	647	rVB2	13775	28734	1.09%	0.189%
10	6.942	651	656	661	rVV3	49770	79543	3.01%	0.524%
11	7.001	661	666	674	rVV2	52809	90773	3.43%	0.598%
12	7.077	674	679	684	rVB3	24726	37089	1.40%	0.244%
13	7.248	701	708	709	rBV2	34842	63102	2.38%	0.416%
14	7.277	709	713	718	rVB2	122330	185382	7.01%	1.222%
15	7.371	726	729	732	rBV5	5641	7149	0.27%	0.047%
16	7.500	745	751	757	rVB2	73815	128942	4.87%	0.850%
17	7.847	804	810	816	rBV	279595	449616	16.99%	2.963%
18	7.911	818	821	825	rVV3	42967	52312	1.98%	0.345%
19	7.970	825	831	837	rVV3	39166	69572	2.63%	0.459%
20	8.047	841	844	847	rBV	5628	8824	0.33%	0.058%
21	8.111	853	855	859	rVB2	8103	9042	0.34%	0.060%
22	8.223	870	874	877	rBV3	20694	23267	0.88%	0.153%
23	8.270	877	882	887	rVV2	172427	283185	10.70%	1.867%
24	8.323	887	891	900	rVB2	54336	83476	3.15%	0.550%
25	8.399	900	904	909	rVB4	9431	16088	0.61%	0.106%
26	8.452	909	913	914	rBV2	6587	7166	0.27%	0.047%
27	8.487	914	919	925	rVB5	21159	35685	1.35%	0.235%
28	8.646	941	946	954	rBV7	16788	32825	1.24%	0.216%
29	8.805	968	973	975	rBV4	13963	23518	0.89%	0.155%
30	8.846	976	980	987	rVB5	16836	37901	1.43%	0.250%
31	8.957	992	999	1004	rVB4	18265	30881	1.17%	0.204%
32	9.151	1026	1032	1035	rBV6	18835	37762	1.43%	0.249%
33	9.275	1050	1053	1057	rBV4	6300	8882	0.34%	0.059%
34	9.433	1075	1080	1084	rVB5	34390	52729	1.99%	0.348%
35	9.539	1092	1098	1099	rBV4	13435	21306	0.81%	0.140%
36	9.639	1112	1115	1120	rVB3	17891	26652	1.01%	0.176%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

37	9.709	1120	1127	1130	rBV4	13148	25479	0.96%	0.168%
38	9.780	1135	1139	1145	rBV4	19293	30742	1.16%	0.203%
39	10.027	1179	1181	1182	rVV	8965	7115	0.27%	0.047%
40	10.044	1182	1184	1187	rVV3	13304	15805	0.60%	0.104%
41	10.074	1187	1189	1194	rVV4	9200	10941	0.41%	0.072%
42	10.132	1197	1199	1202	rBV2	4663	7212	0.27%	0.048%
43	10.273	1219	1223	1226	rVB4	8253	11785	0.45%	0.078%
44	10.638	1278	1285	1292	rBV	406358	737574	27.87%	4.861%
45	10.767	1300	1307	1315	rVB7	31807	62371	2.36%	0.411%
46	10.914	1326	1332	1336	rBV3	26066	53446	2.02%	0.352%
47	11.014	1345	1349	1356	rVB2	31997	56954	2.15%	0.375%
48	11.143	1366	1371	1379	rVB5	18244	39408	1.49%	0.260%
49	11.590	1445	1447	1452	rVB3	9237	13777	0.52%	0.091%
50	11.666	1456	1460	1465	rBV4	10188	17772	0.67%	0.117%
51	11.883	1492	1497	1498	rBV4	6382	9256	0.35%	0.061%
52	12.054	1524	1526	1530	rVB4	9002	9829	0.37%	0.065%
53	12.107	1530	1535	1539	rVB6	8062	12670	0.48%	0.084%
54	12.177	1543	1547	1548	rBV3	18022	16891	0.64%	0.111%
55	12.242	1556	1558	1562	rBV2	4561	8165	0.31%	0.054%
56	12.300	1562	1568	1576	rVB4	26293	50887	1.92%	0.335%
57	12.406	1582	1586	1590	rVB4	7878	11619	0.44%	0.077%
58	12.477	1593	1598	1601	rBV6	17344	30383	1.15%	0.200%
59	12.518	1604	1605	1612	rVB2	17597	19272	0.73%	0.127%
60	12.653	1625	1628	1630	rBV3	7204	7088	0.27%	0.047%
61	12.729	1638	1641	1647	rVV7	9858	17822	0.67%	0.117%
62	12.823	1653	1657	1661	rBV6	8433	11277	0.43%	0.074%
63	12.894	1665	1669	1673	rVB4	14360	23533	0.89%	0.155%
64	13.000	1682	1687	1692	rBV6	10211	12937	0.49%	0.085%
65	13.047	1692	1695	1702	rBV4	28966	41238	1.56%	0.272%
66	13.229	1722	1726	1730	rVB2	35247	48921	1.85%	0.322%
67	13.358	1747	1748	1755	rVB4	11497	20716	0.78%	0.137%
68	13.440	1758	1762	1767	rVV2	31203	51396	1.94%	0.339%
69	13.534	1771	1778	1785	rVB6	36502	77285	2.92%	0.509%
70	13.669	1794	1801	1808	rVV5	22924	53730	2.03%	0.354%
71	13.746	1808	1814	1818	rVV2	55802	90183	3.41%	0.594%
72	13.799	1822	1823	1825	rVV2	13728	13005	0.49%	0.086%
73	13.834	1827	1829	1831	rVV3	16364	16995	0.64%	0.112%
74	13.852	1831	1832	1835	rVV2	13018	11192	0.42%	0.074%
75	13.893	1835	1839	1843	rVB2	12993	20395	0.77%	0.134%
76	13.987	1853	1855	1858	rVB3	9643	9445	0.36%	0.062%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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

77	14.181	1883	1888	1890	rBV2	13339	21934	0.83%	0.145%
78	14.210	1890	1893	1899	rVB2	46562	73337	2.77%	0.483%
79	14.286	1904	1906	1911	rVB3	13207	18841	0.71%	0.124%
80	14.486	1933	1940	1947	rVV	684941	1107959	41.87%	7.303%
81	15.021	2028	2031	2033	rBV4	7768	9659	0.36%	0.064%
82	15.115	2041	2047	2059	rVB	431060	703107	26.57%	4.634%
83	15.350	2083	2087	2090	rBV4	4983	8381	0.32%	0.055%
84	15.485	2104	2110	2114	rVB2	18156	31475	1.19%	0.207%
85	15.620	2128	2133	2143	rBV3	33901	71980	2.72%	0.474%
86	16.748	2321	2325	2329	rVB3	7988	11864	0.45%	0.078%
87	16.783	2329	2331	2335	rBV3	7071	8899	0.34%	0.059%
88	16.889	2342	2349	2363	rBV	1373692	2206151	83.37%	14.541%
89	17.236	2402	2408	2417	rVV	1119255	1685046	63.67%	11.106%
90	17.324	2421	2423	2430	rVB3	26098	45571	1.72%	0.300%
91	17.970	2528	2533	2535	rVB2	5860	9798	0.37%	0.065%
92	18.834	2678	2680	2683	rBV3	7785	7175	0.27%	0.047%
93	19.034	2711	2714	2717	rBV4	5530	7467	0.28%	0.049%
94	19.410	2775	2778	2780	rBV3	6244	6928	0.26%	0.046%
95	19.633	2809	2816	2823	rBV2	39979	70622	2.67%	0.465%
96	20.156	2903	2905	2908	rBV3	6093	7963	0.30%	0.052%
97	20.591	2977	2979	2980	rBV	10984	8538	0.32%	0.056%
98	21.484	3125	3131	3140	rBV	1648113	2559969	96.74%	16.873%
99	23.017	3390	3392	3393	rBV2	12998	8028	0.30%	0.053%
100	24.527	3638	3649	3662	rBV	956741	2646340	100.00%	17.442%

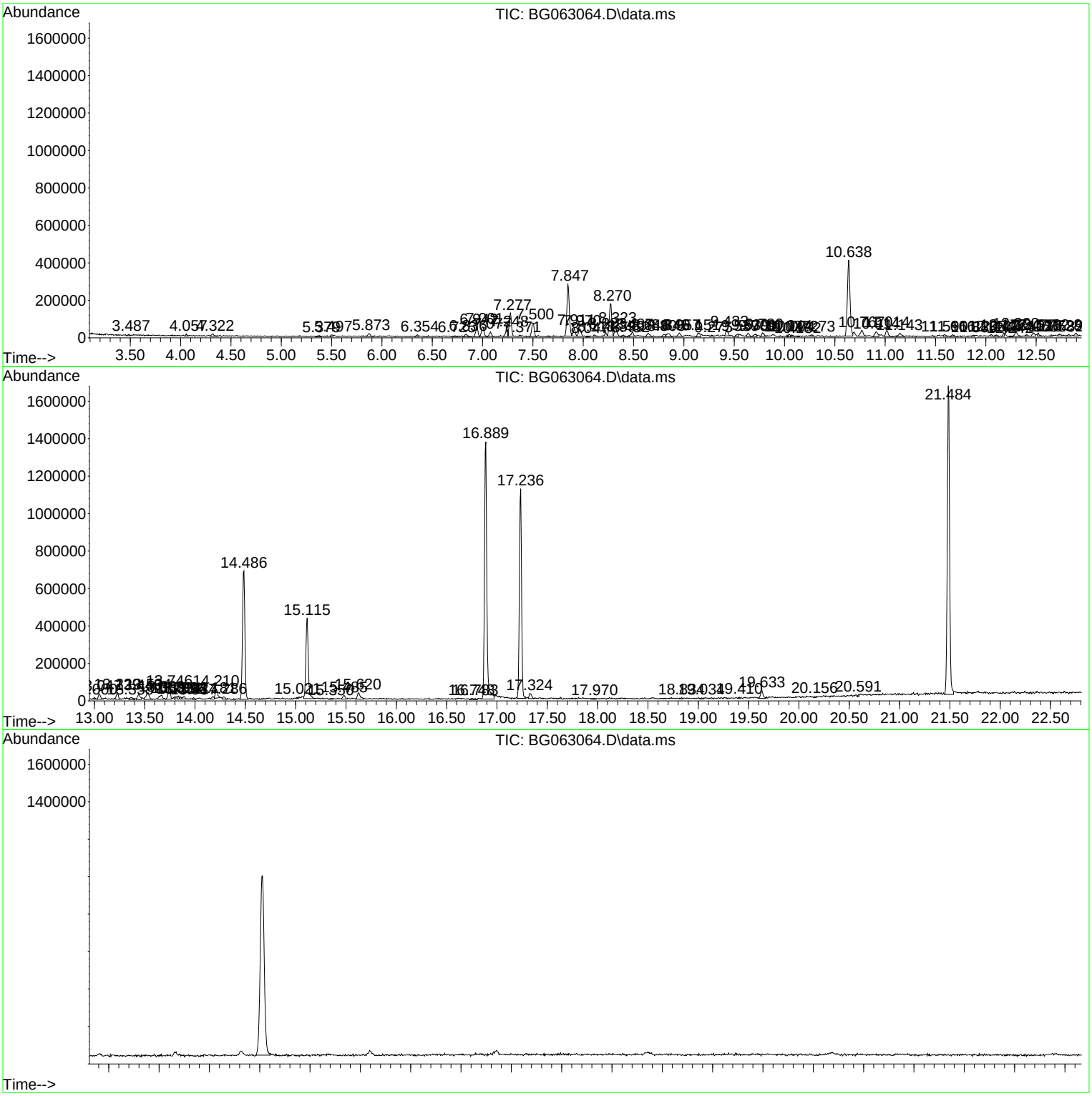
Sum of corrected areas: 15171867

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

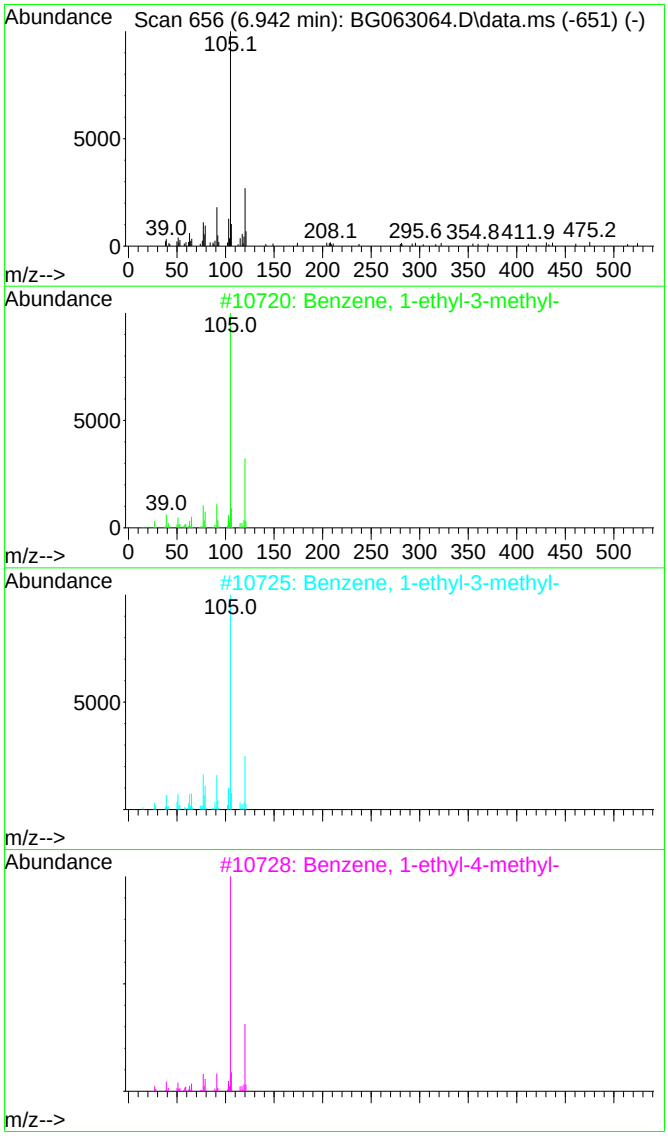
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

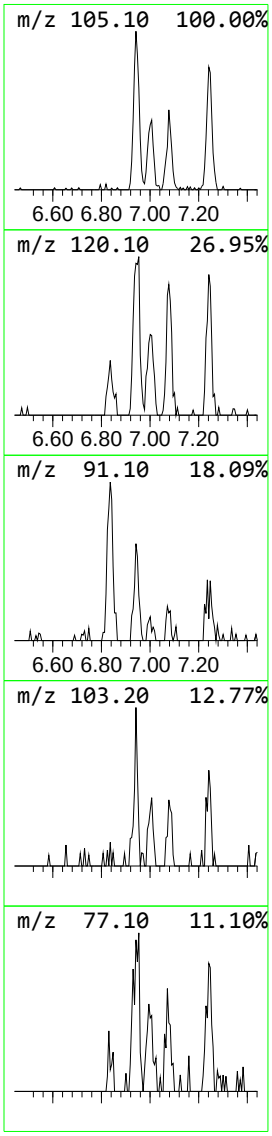
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.942	3.54 ng/ul	79543	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	68
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	68
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	64
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	64
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

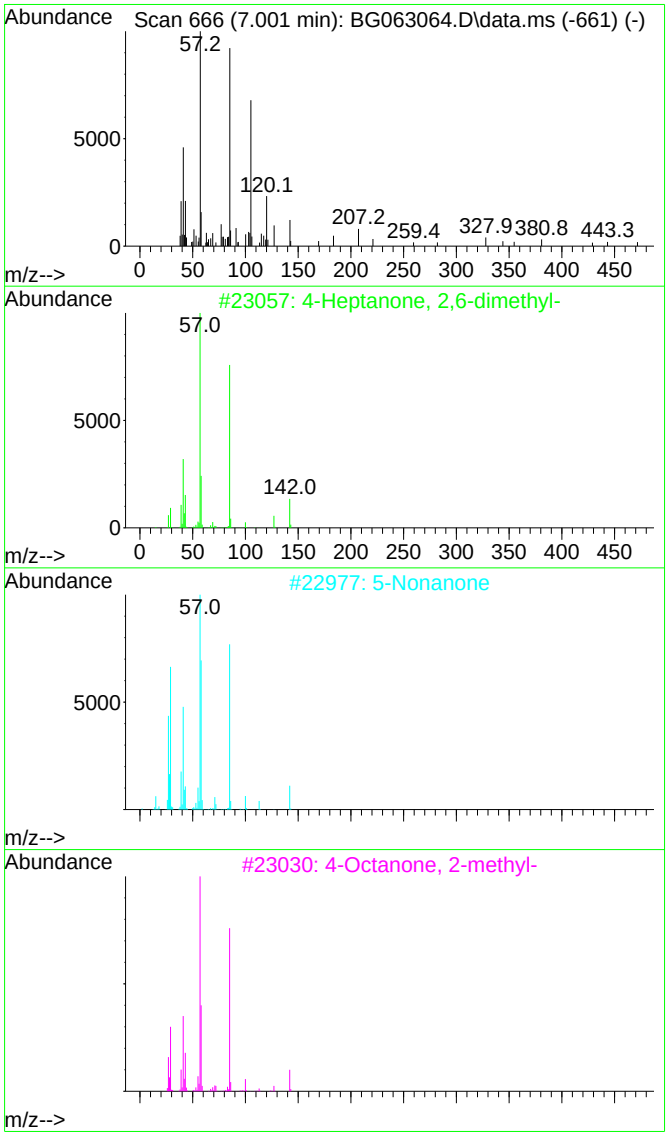
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

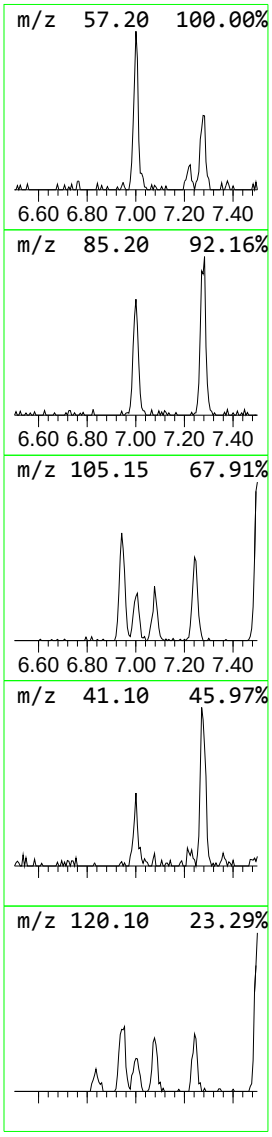
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 4-Heptanone, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.001	4.04 ng/ul	90773	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	59
2	5-Nonanone	142	C9H18O	000502-56-7	58
3	4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	58
4	4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	53
5	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

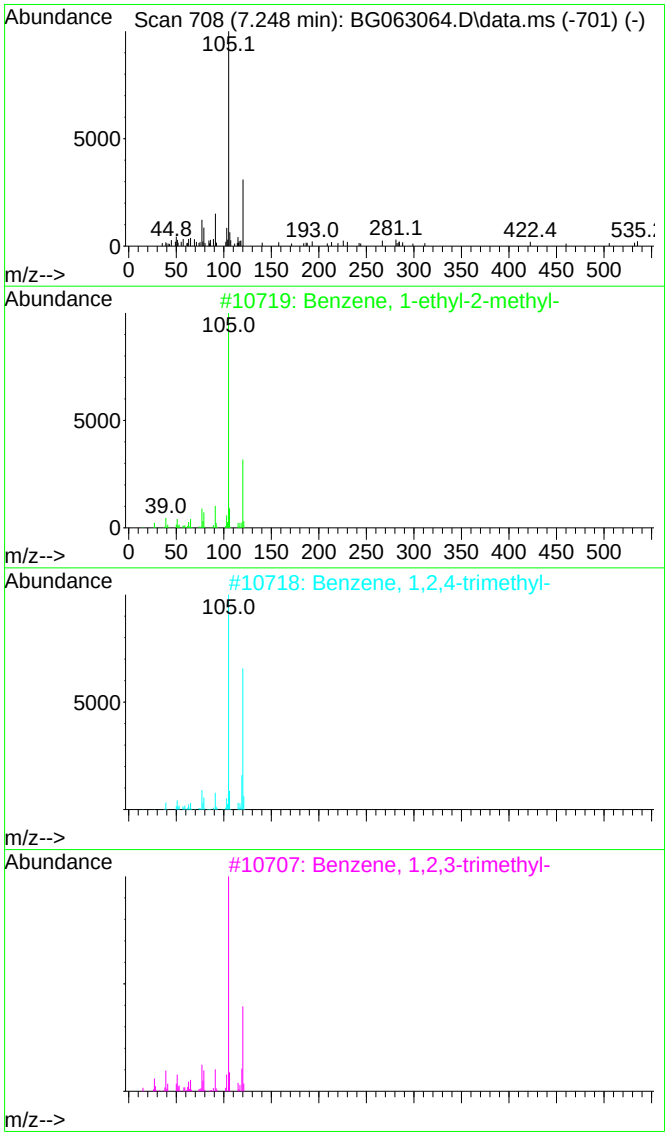
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

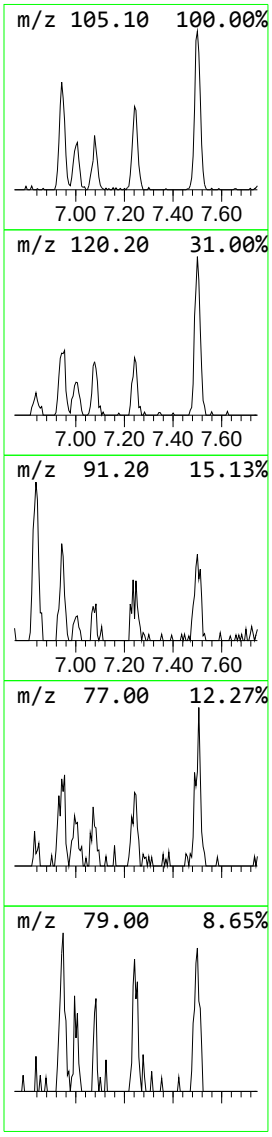
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.248	2.81 ng/ul	63102	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	58
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	58
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	58
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

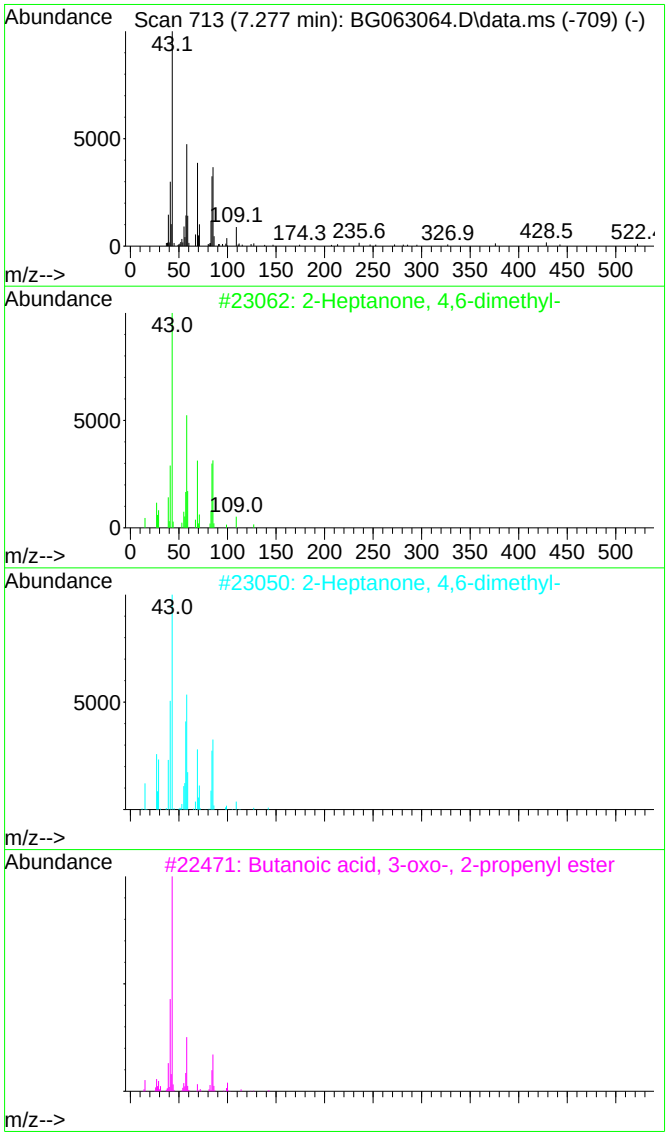
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

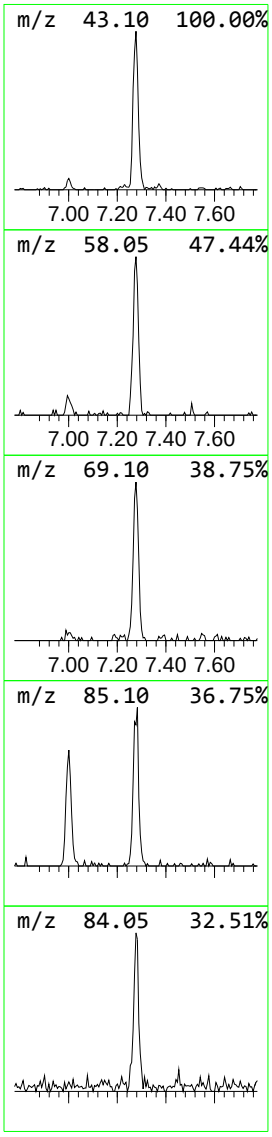
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Heptanone, 4,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.277	8.25 ng/ul	185382	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	72
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	56
3	Butanoic acid, 3-oxo-, 2-propeny...	142	C7H10O3	001118-84-9	53
4	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	43
5	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

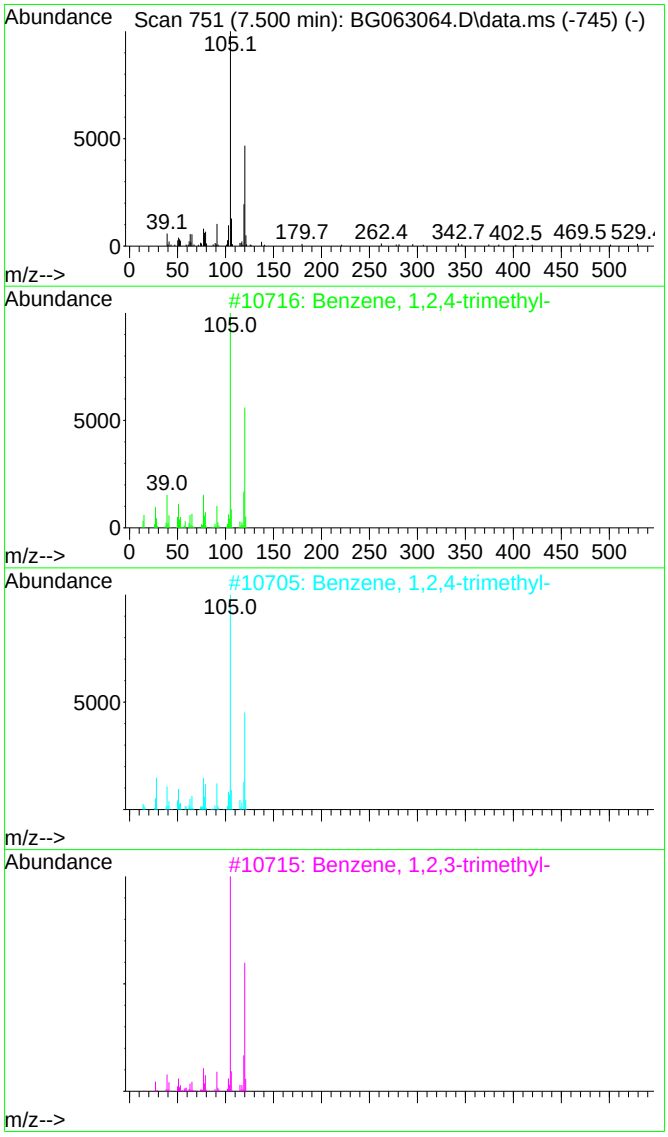
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

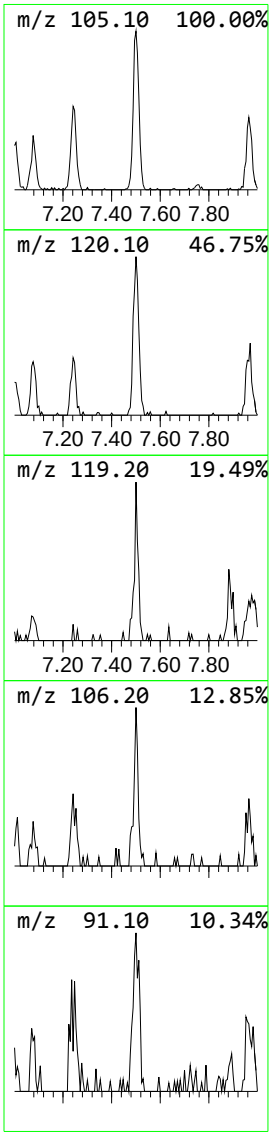
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.500	5.74 ng/ul	128942	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	83
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	80
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	74
4	Mesitylene	120	C9H12	000108-67-8	74
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	72





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Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

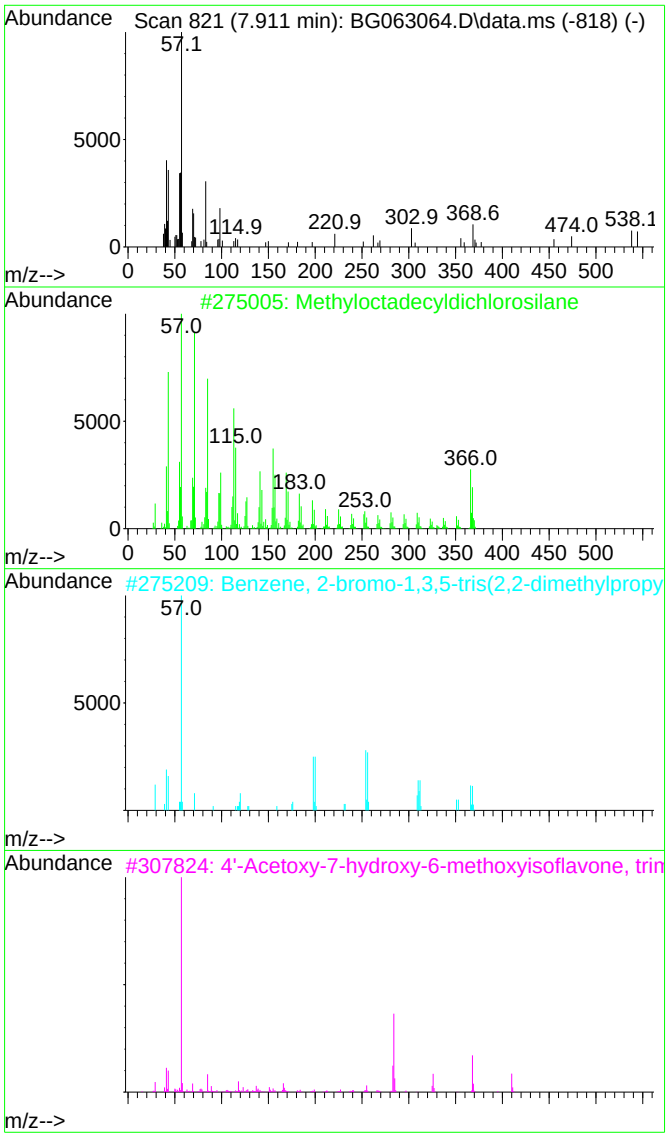
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

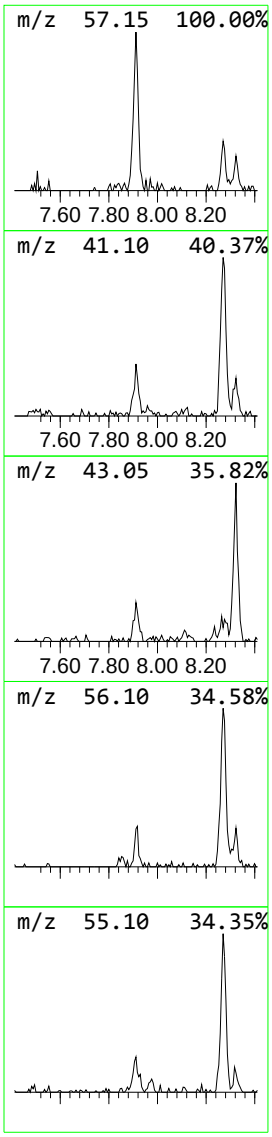
****

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.911	2.33 ng/ul	52312	1,4-Dichlorobenzene-d4	7.853

Hit# of 4	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyloctadecyldichlorosilane	366	C19H40Cl2Si	005157-75-5	1
2	Benzene, 2-bromo-1,3,5-tris(2,2-...	366	C21H35Br	025347-03-9	1
3	4'-Acetoxy-7-hydroxy-6-methoxyis...	410	C23H22O7	1000449-16-8	1
4	Cholest-5-en-3-ol (3.beta.)-, pr...	442	C30H50O2	000633-31-8	1





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

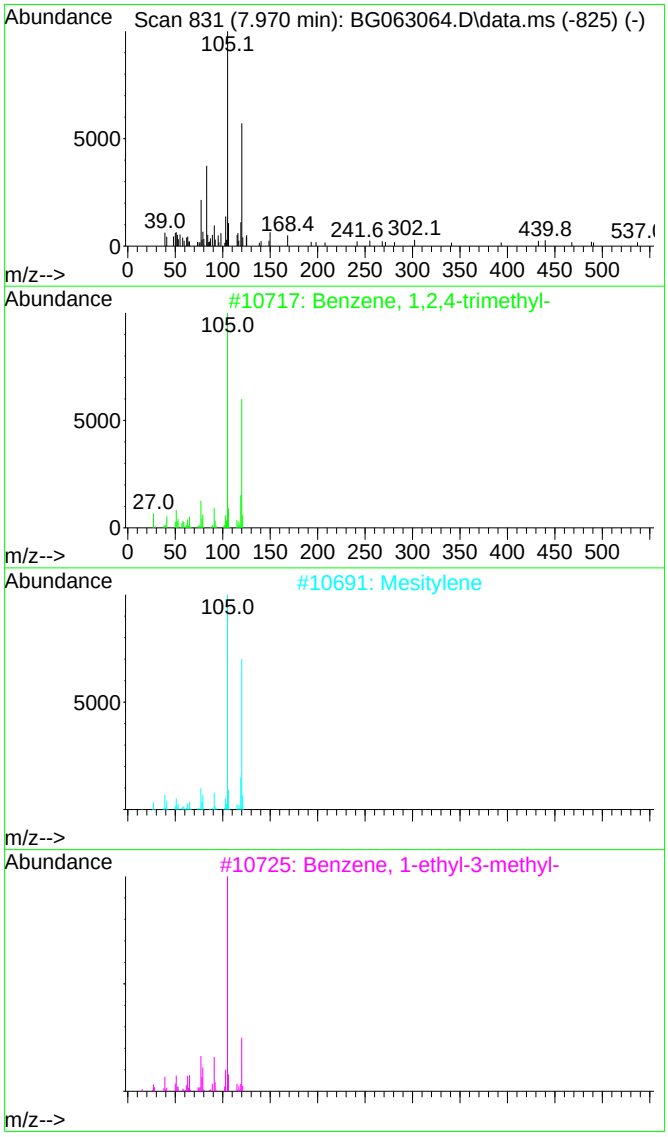
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

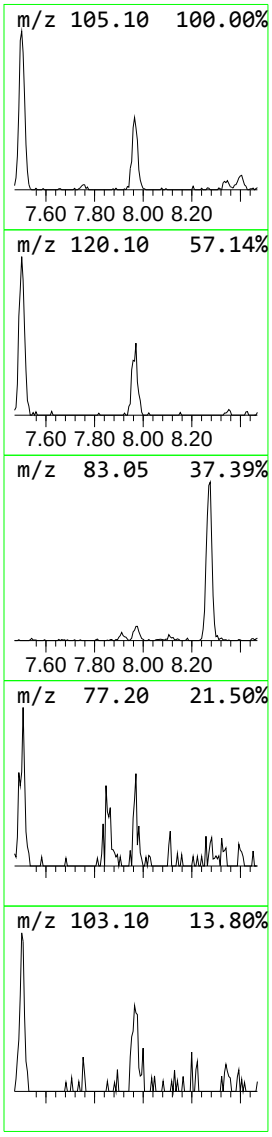
****

Peak Number 7 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.970	3.09 ng/ul	69572	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	60
2	Mesitylene	120	C9H12	000108-67-8	53
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	53
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	53
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	49





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Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

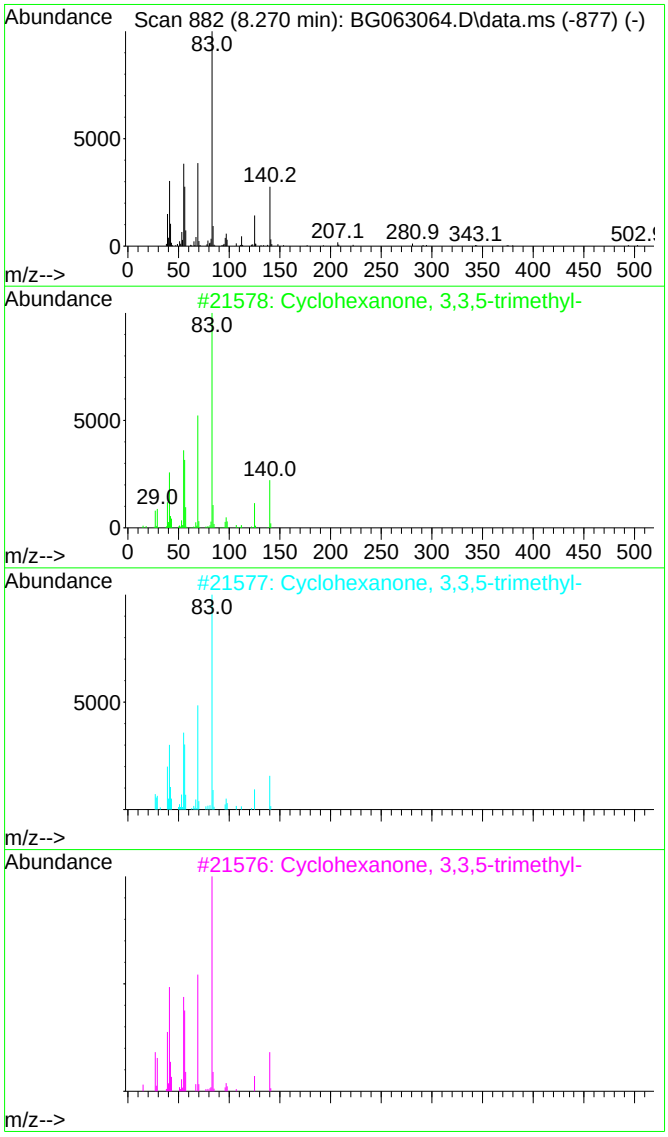
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

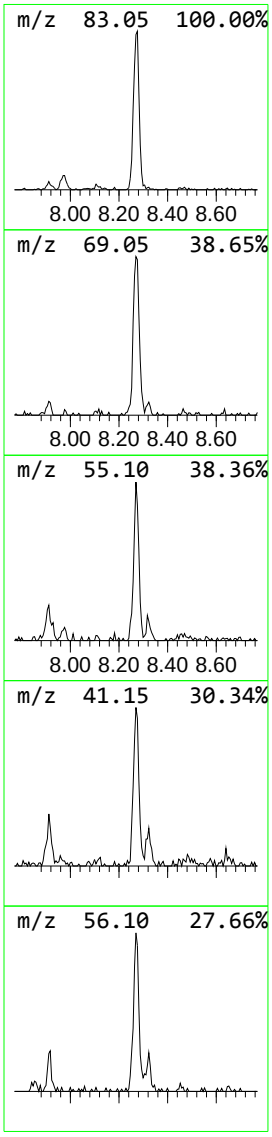
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclohexanone, 3,3,5-trimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.270	12.60 ng/ul	283185	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
5	Propanenitrile, 3-[(2-methylprop...	126	C7H14N2	014278-96-7	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

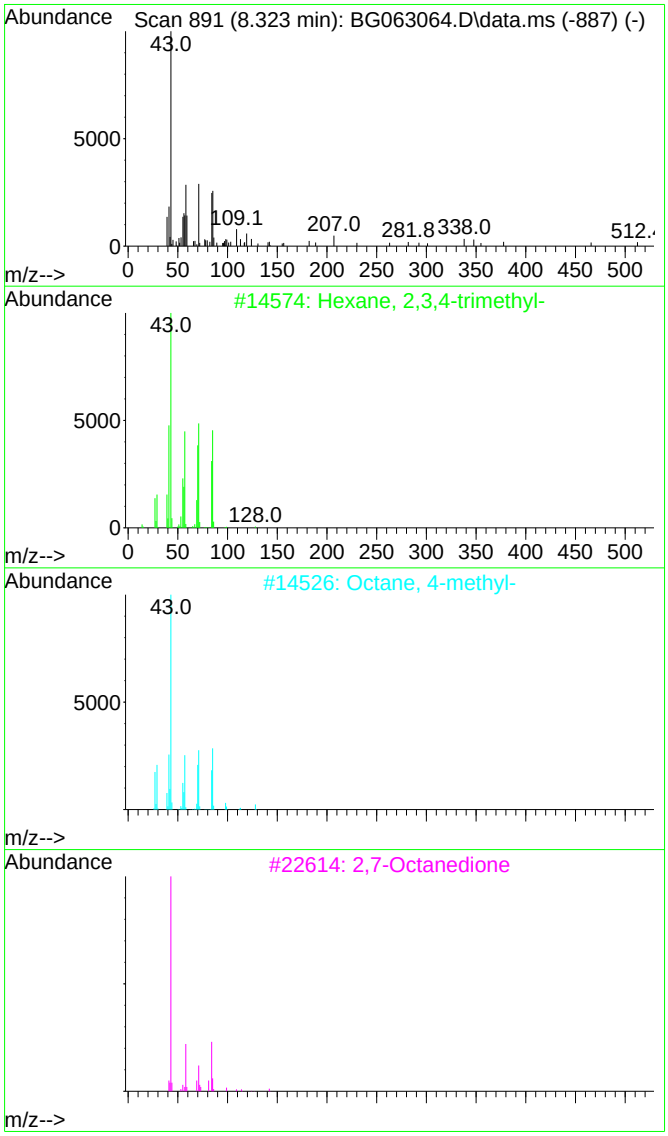
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

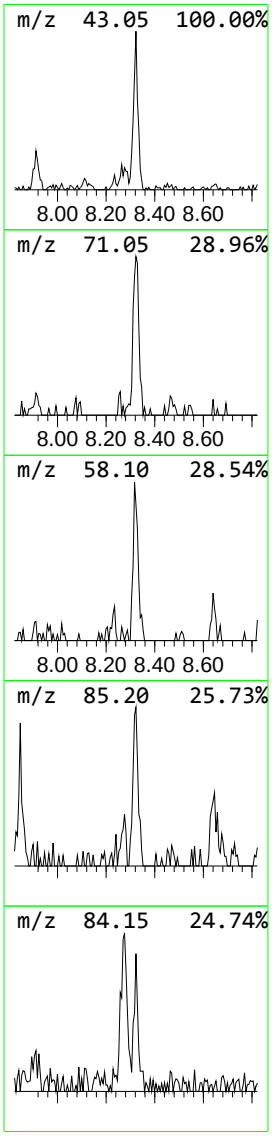
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.323	3.71 ng/ul	83476	1,4-Dichlorobenzene-d4	7.853

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	43
2	Octane, 4-methyl-	128	C9H20	002216-34-4	43
3	2,7-Octanedione	142	C8H14O2	001626-09-1	43
4	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	42
5	Octane, 4-methyl-	128	C9H20	002216-34-4	37





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063064.D
Acq On : 26 Sep 2024 14:26
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

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

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-ethy...	6.942	3.5	ng/ul	79543	1	7.853	449616	20.0
4-Heptanone, 2,...	7.001	4.0	ng/ul	90773	1	7.853	449616	20.0
Benzene, 1-ethy...	7.248	2.8	ng/ul	63102	1	7.853	449616	20.0
2-Heptanone, 4,...	7.277	8.3	ng/ul	185382	1	7.853	449616	20.0
Benzene, 1,2,4-...	7.500	5.7	ng/ul	128942	1	7.853	449616	20.0
unknown-01	7.911	2.3	ng/ul	52312	1	7.853	449616	20.0
Mesitylene	7.970	3.1	ng/ul	69572	1	7.853	449616	20.0
Cyclohexanone, ...	8.270	12.6	ng/ul	283185	1	7.853	449616	20.0
(DEL) Alkane: S...	8.323	3.7	ng/ul	83476	1	7.853	449616	20.0