LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

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-BG090924.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062783.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	35	42	75	rVB	3290411	6679837	100.00%	22.367%
2	3.768	112	116	130	rBV	55841	102464	1.53%	0.343%
3	4.103	169	173	179	rBV	20604	31212	0.47%	0.105%
4	4.174	182	185	199	rVB4	17584	43168	0.65%	0.145%
5	4.632	257	263	284	rVB	377777	676347	10.13%	2.265%
6	5.020	324	329	333	rBV3	4776	8090	0.12%	0.027%
7	5.196	354	359	366	rBV	22780	38553	0.58%	0.129%
8	5.419	393	397	402	rBV	10449	15023	0.22%	0.050%
9	5.555	413	420	432	rBV	92451	205843	3.08%	0.689%
10	5.919	474	482	497	rBV	210665	347163	5.20%	1.162%
11	6.882	639	646	658	rBV	199119	345593	5.17%	1.157%
12	6.994	658	665	670	rVV2	7992	13143	0.20%	0.044%
13	7.065	670	677	684	rVV2	76628	139016	2.08%	0.465%
14	7.123	684	687	695	rVV	12431	19761	0.30%	0.066%
15	7.229	697	705	712	rBV	174037	285922	4.28%	0.957%
16	7.294	712	716	728	rVB	37345	61723	0.92%	0.207%
17	7.435	733	740	750	rVB	204070	350935	5.25%	1.175%
18	7.552	754	760	767	rVB	32562	50437	0.76%	0.169%
19	7.899	813	819	830	rBV	160264	284941	4.27%	0.954%
20	8.016	833	839	849	rVB	135847	235516	3.53%	0.789%
21	8.210	864	872	874	rBV3	7662	11410	0.17%	0.038%
22	8.251	874	879	886	rVB2	30226	53755	0.80%	0.180%
23	8.392	898	903	907	rBV5	7415	11147	0.17%	0.037%
24	8.451	909	913	920	rVB3	7217	11240	0.17%	0.038%
25	8.539	920	928	932	rBV3	17766	30985	0.46%	0.104%
26	8.604	932	939	948	rVV	193261	338576	5.07%	1.134%
27	8.704	951	956	965	rVB	18575	34503	0.52%	0.116%
28	8.904	986	990	997	rVB3	5657	8485	0.13%	0.028%
29	9.004	1002	1007	1009	rVV2	17360	25919	0.39%	0.087%
30	9.051	1009	1015	1031	rVB	222280	422489	6.32%	1.415%
31	9.250	1044	1049	1057	rBV6	4844	12326	0.18%	0.041%
32	9.338	1057	1064	1070	rVV3	14447	28451	0.43%	0.095%
33	9.773	1131	1138	1148	rBV	183179	323316	4.84%	1.083%
34	9.885	1152	1157	1166	rVB7	5674	12254	0.18%	0.041%
35	10.102	1187	1194	1200	rVB3	8398	14137	0.21%	0.047%
36	10.314	1223	1230	1239	rBV	289072	510995	7.65%	1.711%

LSC Area Percent Report

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Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

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-BG090924.MA.M
Title : SVOA CALIBRATION

37	10.390	1239	1243	1250	rVV4	5190	10792	0.16%	0.036%
38	10.608	1276	1280	1287	rVB3	11386	16245	0.24%	0.054%
39	10.690	1287	1294	1300	rBV	214540	395699	5.92%	1.325%
40	10.743	1300	1303	1308	rVV2	29878	42691	0.64%	0.143%
41	10.825	1308	1317	1329	rVV	169469	311324	4.66%	1.042%
42	11.177	1373	1377	1383	rBV4	6957	13972	0.21%	0.047%
43	11.360	1403	1408	1416	rVB6	6572	13501	0.20%	0.045%
44	11.501	1427	1432	1437	rBV4	10264	15291	0.23%	0.051%
45	11.553	1437	1441	1445	rVV2	8530	14001	0.21%	0.047%
46	11.800	1478	1483	1487	rBV4	4351	8083	0.12%	0.027%
47	11.918	1498	1503	1512	rBV4	11949	22899	0.34%	0.077%
48	12.047	1520	1525	1532	rBV8	7171	15911	0.24%	0.053%
49	12.247	1554	1559	1562	rBV4	5407	10259	0.15%	0.034%
50	12.306	1566	1569	1578	rVV6	7335	14427	0.22%	0.048%
51	12.405	1581	1586	1591	rBV3	5628	9307	0.14%	0.031%
52	12.505	1596	1603	1611	rBV2	19496	30394	0.46%	0.102%
53	12.658	1624	1629	1641	rVB5	18375	39606	0.59%	0.133%
54	12.928	1670	1675	1680	rVV3	20961	40988	0.61%	0.137%
55	12.981	1680	1684	1688	rVV5	4213	8371	0.13%	0.028%
56	13.257	1725	1731	1736	rBV7	10979	23552	0.35%	0.079%
57	13.334	1738	1744	1752	rVB7	8373	20442	0.31%	0.068%
58	13.528	1771	1777	1780	rBV5	9895	17002	0.25%	0.057%
59	13.739	1809	1813	1820	rVB5	9473	15962	0.24%	0.053%
60	13.810	1823	1825	1830	rBV6	4874	8042	0.12%	0.027%
61	13.933	1839	1846	1852	rBV	716825	1023764	15.33%	3.428%
62	14.062	1862	1868	1872	rBV3	29718	39430	0.59%	0.132%
63	14.139	1876	1881	1887	rBV4	14989	26468	0.40%	0.089%
64	14.221	1887	1895	1902	rVV	752044	1168265	17.49%	3.912%
65	14.297	1904	1908	1915	rVV4	7726	13745	0.21%	0.046%
66	14.526	1936	1947	1956	rBV	473993	724944	10.85%	2.427%
67	14.720	1973	1980	1992	rBV2	88170	172029	2.58%	0.576%
68	14.867	2001	2005	2013	rBV6	9191	15142	0.23%	0.051%
69	14.944	2013	2018	2025	rBV4	9348	13361	0.20%	0.045%
70	15.519	2109	2116	2130	rBV	1100802	1616451	24.20%	5.413%
71	15.637	2130	2136	2146	rBV	332426	491728	7.36%	1.647%
72	15.884	2171	2178	2182	rBV	42337	62123	0.93%	0.208%
73	16.195	2227	2231	2241	rBV5	11308	20226	0.30%	0.068%
74	17.141	2389	2392	2396	rVB2	5852	8400	0.13%	0.028%
75	17.270	2407	2414	2420	rBV	684380	994139	14.88%	3.329%
76	17.370	2424	2431	2442	rVB	1334262	1953598	29.25%	6.542%

LSC Area Percent Report

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Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

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-BG090924.MA.M
Title : SVOA CALIBRATION

77	18.163	2561	2566	2573	rBV3	40768	68232	1.02%	0.228%
78	18.645	2645	2648	2653	rBV4	5794	10867	0.16%	0.036%
79	19.227	2742	2747	2753	rVB3	24449	33535	0.50%	0.112%
80	19.297	2754	2759	2761	rVV	21898	33254	0.50%	0.111%
81	19.327	2761	2764	2770	rVB	37532	52450	0.79%	0.176%
82	19.444	2779	2784	2790	rBV6	14064	25544	0.38%	0.086%
83	19.662	2813	2821	2833	rBV	1944323	2783436	41.67%	9.320%
84	20.185	2908	2910	2917	rVB8	6026	11760	0.18%	0.039%
85	20.337	2931	2936	2939	rBV7	5644	10224	0.15%	0.034%
86	21.119	3063	3069	3071	rBV6	6623	12441	0.19%	0.042%
87	21.189	3077	3081	3089	rVB9	20660	32864	0.49%	0.110%
88	21.412	3117	3119	3125	rVB7	10113	15164	0.23%	0.051%
89	21.512	3130	3136	3144	rBV	762892	1273473	19.06%	4.264%
90	21.712	3166	3170	3175	rBV8	12796	18752	0.28%	0.063%
91	22.294	3265	3269	3275	rVB8	14758	22541	0.34%	0.075%
92	22.946	3376	3380	3387	rVB5	15472	28921	0.43%	0.097%
93	23.128	3407	3411	3418	rVB7	20563	36630	0.55%	0.123%
94	23.604	3487	3492	3494	rBV4	16591	30825	0.46%	0.103%
95	24.291	3605	3609	3611	rBV3	8612	13417	0.20%	0.045%
96	24.368	3611	3622	3642	rVB	1017081	2715583	40.65%	9.093%
97	24.574	3647	3657	3672	rVV	488272	1328954	19.90%	4.450%
98	25.655	3834	3841	3842	rBV5	18135	30569	0.46%	0.102%
99	26.260	3942	3944	3950	rVB7	5467	7907	0.12%	0.026%
100	28.005	4235	4241	4242	rBV6	12694	19491	0.29%	0.065%

Sum of corrected areas: 29864083

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
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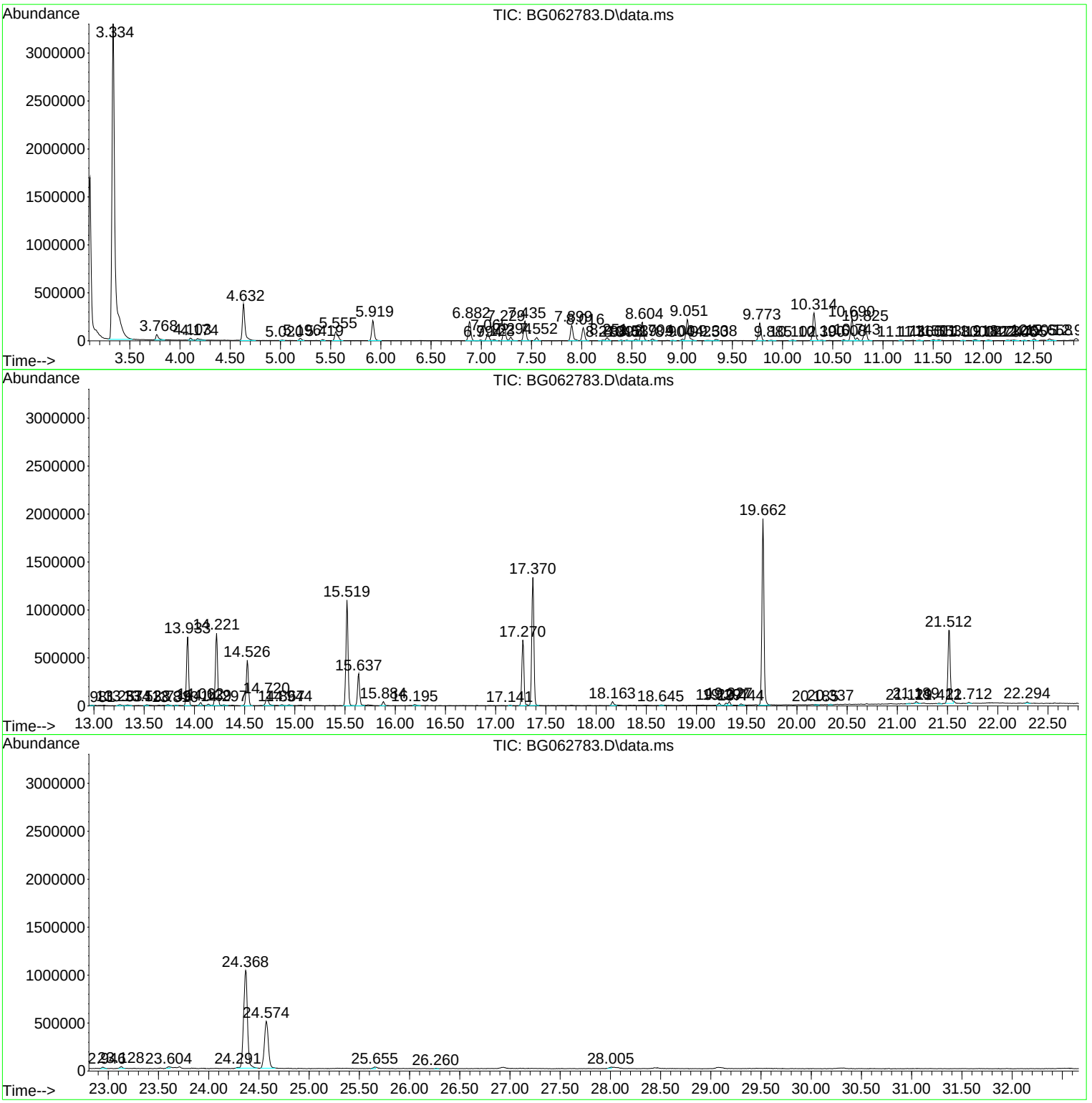
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.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\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

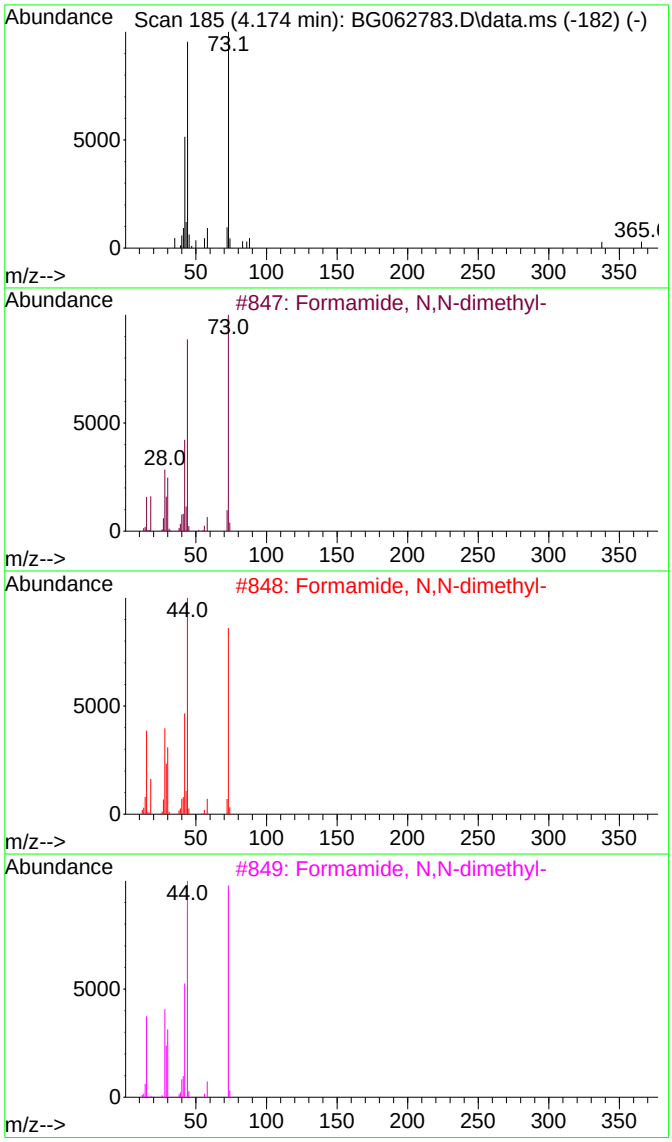
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

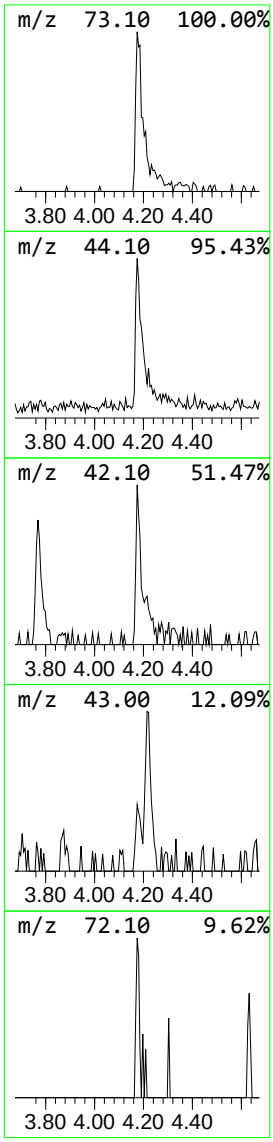
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Formamide, N,N-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.174	3.03 ng/ul	43168	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	86
2	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	86
3	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	72
4	Glycinamide, N(2)-methyl-	88	C3H8N2O	1000452-55-9	64
5	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

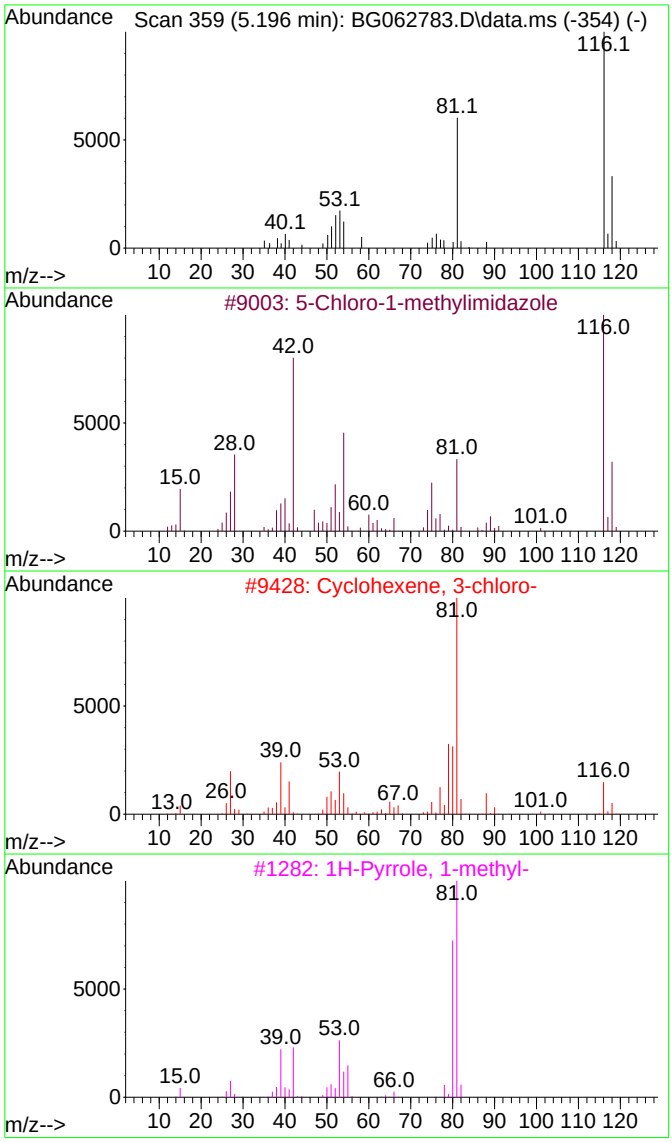
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

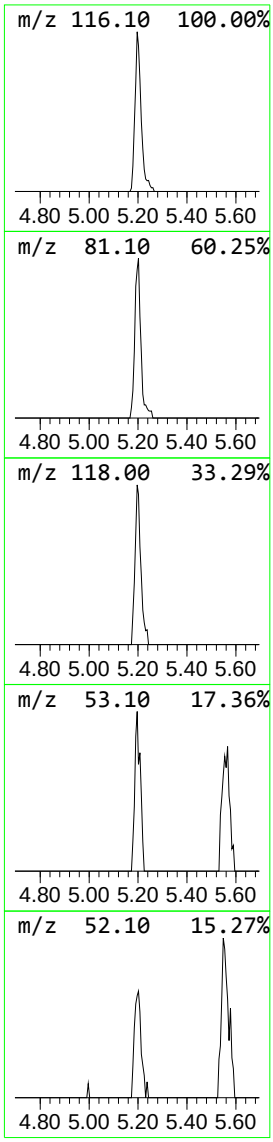
****

Peak Number 5 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.196	2.71 ng/ul	38553	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Chloro-1-methylimidazole	116	C4H5ClN2	000872-49-1	45
2	Cyclohexene, 3-chloro-	116	C6H9Cl	002441-97-6	42
3	1H-Pyrrole, 1-methyl-	81	C5H7N	000096-54-8	11
4	Cyclopropanecarboxaldehyde, meth...	82	C5H6O	142423-24-3	10
5	2H-Thiopyran, tetrahydro-4-methyl-	116	C6H12S	005161-17-1	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

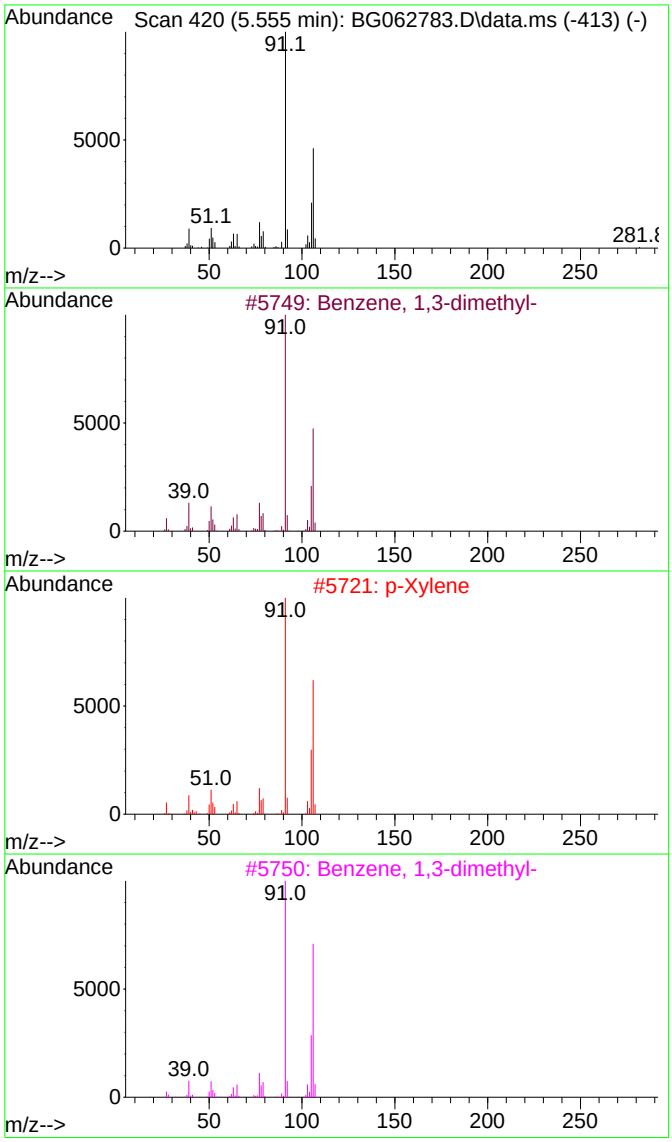
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

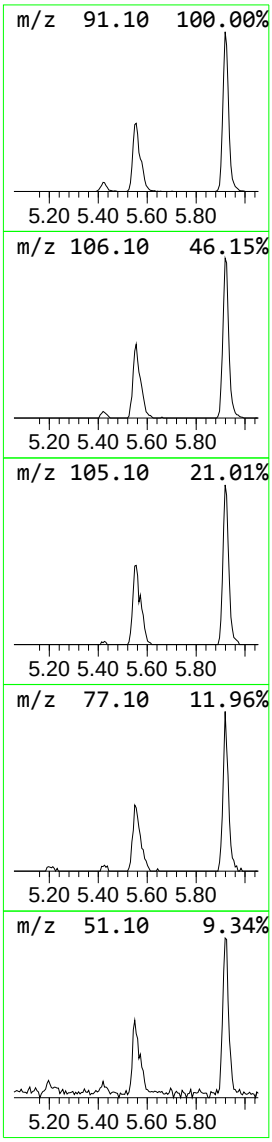
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.555	14.45 ng/ul	205843	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	p-Xylene	106	C8H10	000106-42-3	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4	o-Xylene	106	C8H10	000095-47-6	95
5	p-Xylene	106	C8H10	000106-42-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

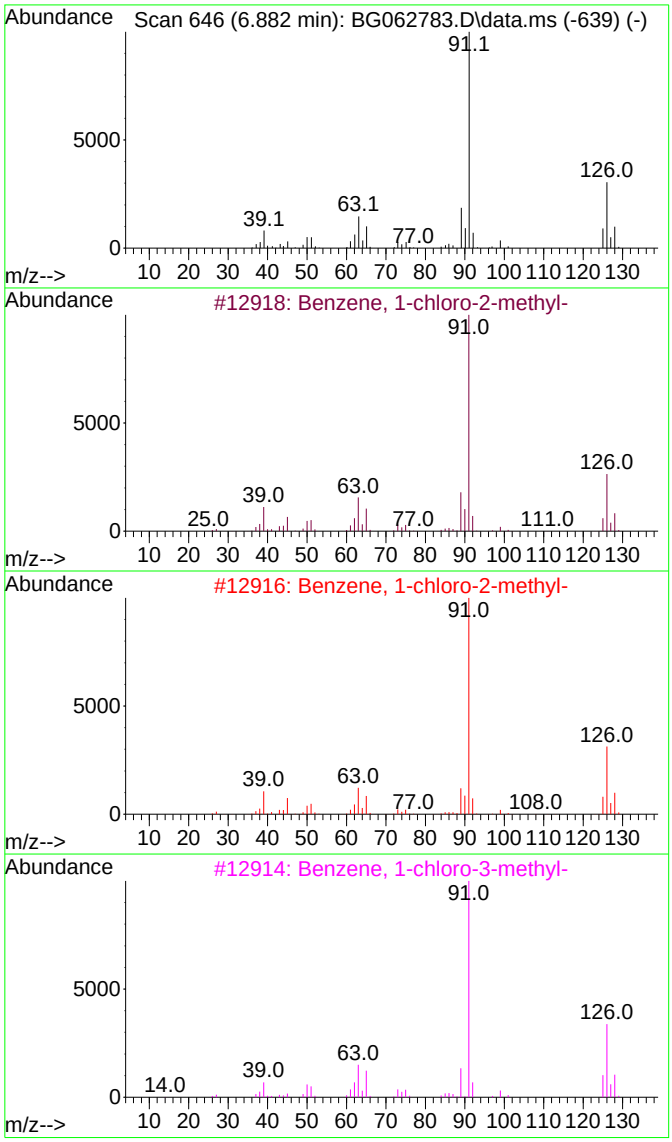
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

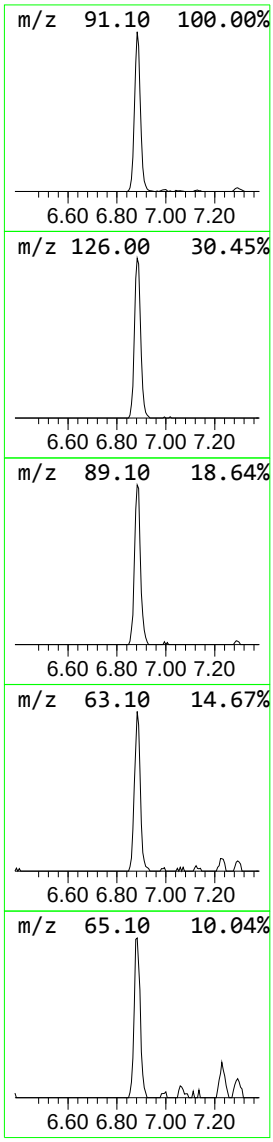
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-chloro-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.882	24.26 ng/ul	345593	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	96
2	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	95
3	Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	95
4	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	95
5	Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

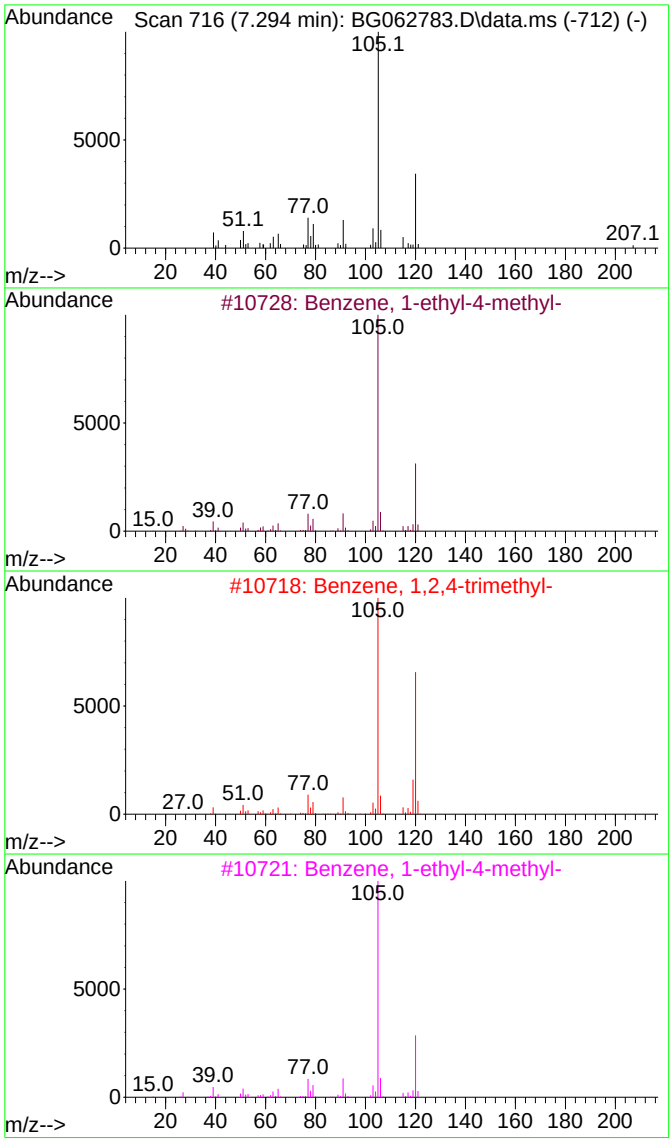
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

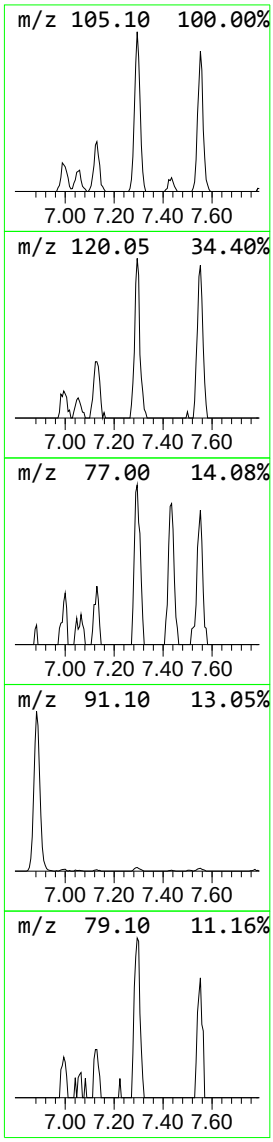
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.294	4.33 ng/ul	61723	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





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Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

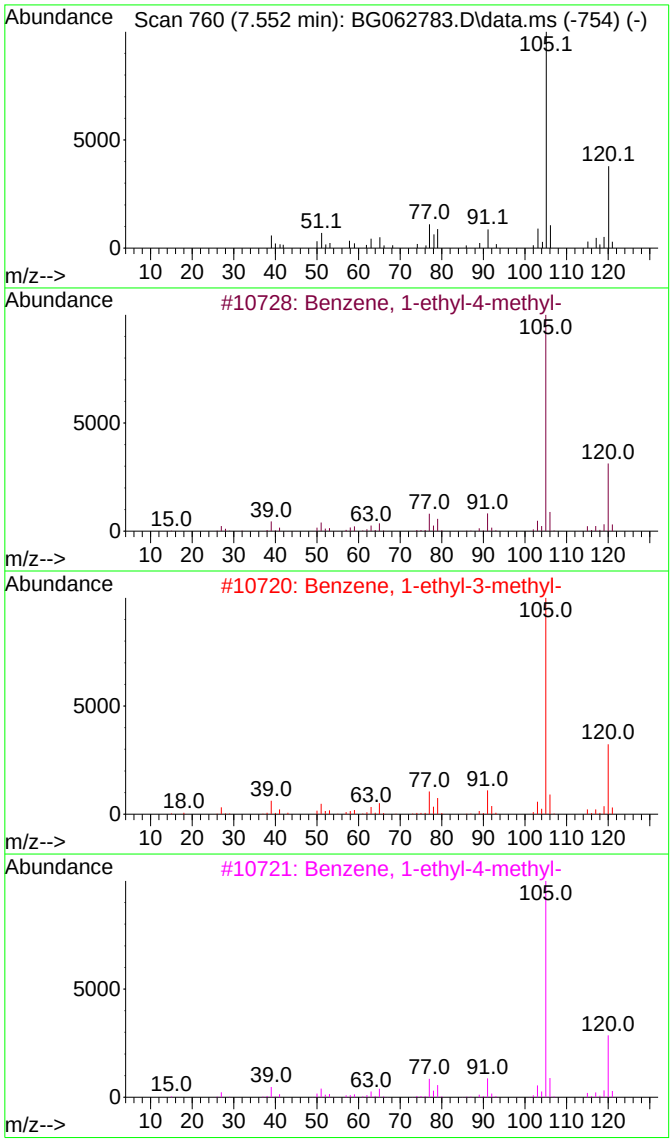
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

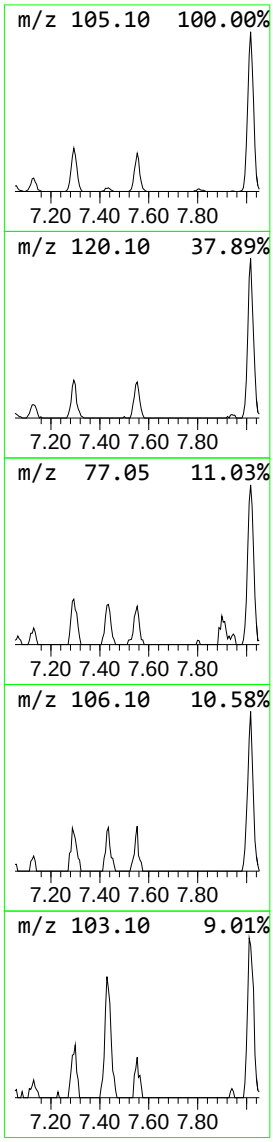
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.552	3.54 ng/ul	50437	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5	Mesitylene	120	C9H12	000108-67-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

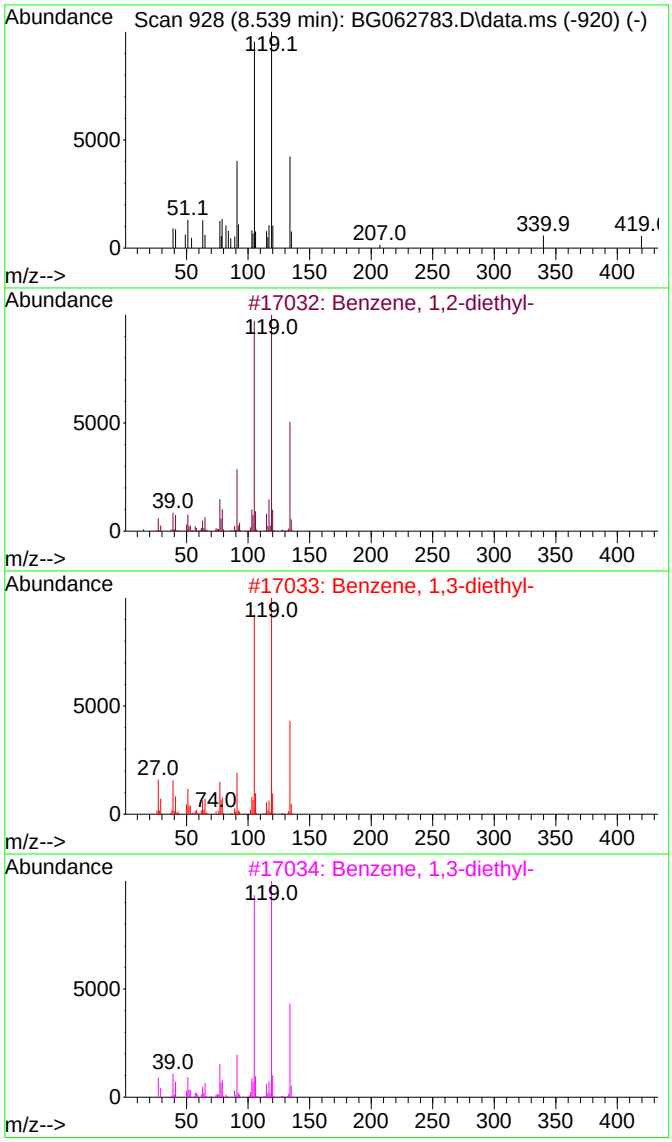
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

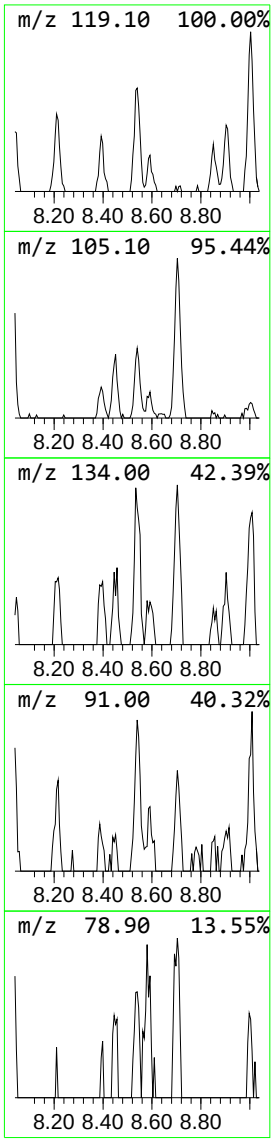
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,2-diethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.539	2.17 ng/ul	30985	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

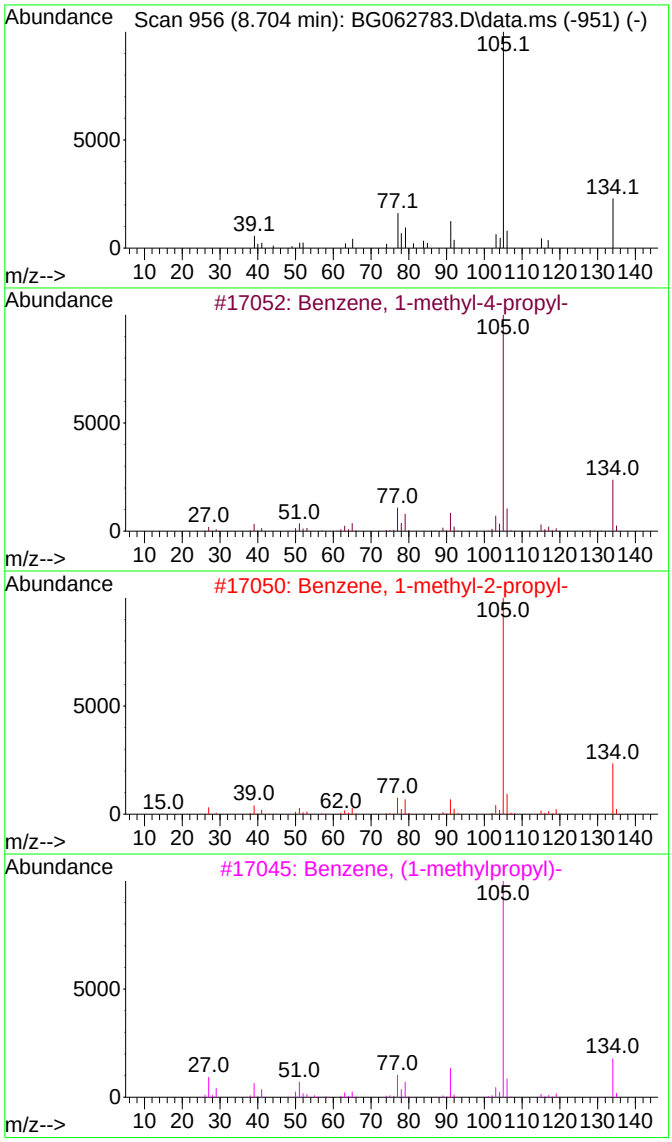
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

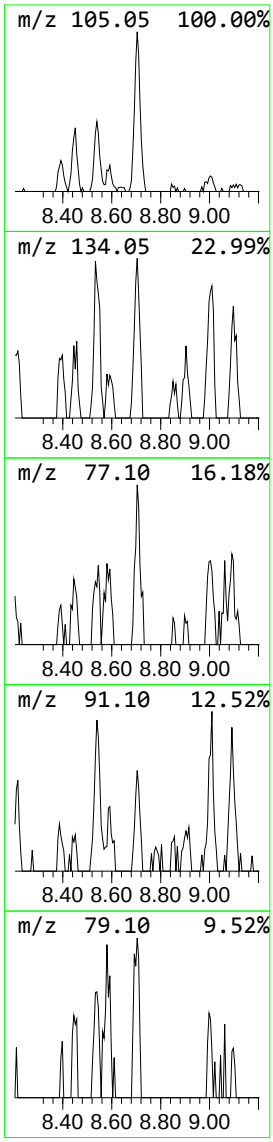
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-methyl-4-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.704	2.42 ng/ul	34503	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	86
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062783.D
Acq On : 13 Sep 2024 11:46
Operator : JU/RC
Sample : P3952-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B38

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.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
Formamide, N,N-...	4.174	3.0	ng/ul	43168	1	7.899	284941	20.0
unknown-01	5.196	2.7	ng/ul	38553	1	7.899	284941	20.0
Benzene, 1,3-di...	5.555	14.4	ng/ul	205843	1	7.899	284941	20.0
Benzene, 1-chlo...	6.882	24.3	ng/ul	345593	1	7.899	284941	20.0
Benzene, 1-ethy...	7.294	4.3	ng/ul	61723	1	7.899	284941	20.0
Benzene, 1-ethy...	7.552	3.5	ng/ul	50437	1	7.899	284941	20.0
Benzene, 1,2-di...	8.539	2.2	ng/ul	30985	1	7.899	284941	20.0
Benzene, 1-meth...	8.704	2.4	ng/ul	34503	1	7.899	284941	20.0