LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8Y9DL

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_P\Methods\SFAM-EPA-BP101122.M
Title : SVOA CALIBRATION

Signal : TIC: BP012124.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	12	15	25	rVB	115045	171727	4.74%	0.396%
2	3.311	51	55	66	rVB	222468	296889	8.20%	0.684%
3	3.705	120	122	135	rBV	217281	341486	9.43%	0.787%
4	3.917	156	158	163	rVB2	10705	14006	0.39%	0.032%
5	4.017	172	175	179	rBV	19015	22371	0.62%	0.052%
6	4.469	246	252	258	rVB	183069	263905	7.29%	0.608%
7	5.152	362	368	376	rVB	178020	263662	7.28%	0.608%
8	5.417	410	413	420	rVB8	7926	16147	0.45%	0.037%
9	5.593	440	443	456	rVB2	22008	42060	1.16%	0.097%
10	5.781	469	475	480	rBV	14034	23433	0.65%	0.054%
11	6.128	524	534	539	rVB	11689	24256	0.67%	0.056%
12	6.440	582	587	593	rBV5	12168	27221	0.75%	0.063%
13	6.516	597	600	606	rVB3	11616	18063	0.50%	0.042%
14	7.016	678	685	698	rBV	113098	228611	6.32%	0.527%
15	7.181	705	713	731	rBV	397228	667772	18.45%	1.539%
16	7.381	740	747	755	rBV	378007	624816	17.26%	1.440%
17	7.505	764	768	775	rVB8	12570	25353	0.70%	0.058%
18	7.846	819	826	836	rBV	881938	1430259	39.52%	3.297%
19	7.934	836	841	851	rVB4	29551	76006	2.10%	0.175%
20	8.216	883	889	897	rVB	109895	187499	5.18%	0.432%
21	8.334	902	909	915	rVB6	10604	18081	0.50%	0.042%
22	8.510	928	939	942	rBV4	18799	38086	1.05%	0.088%
23	8.563	942	948	958	rVV	324458	548647	15.16%	1.265%
24	8.646	958	962	966	rVB7	8936	14261	0.39%	0.033%
25	8.763	977	982	989	rVB8	8450	20004	0.55%	0.046%
26	8.840	989	995	1005	rBV	144638	248702	6.87%	0.573%
27	8.957	1009	1015	1019	rVV3	43754	78239	2.16%	0.180%
28	9.016	1019	1025	1035	rVV	422993	735173	20.31%	1.695%
29	9.216	1052	1059	1068	rVB	13029	33809	0.93%	0.078%
30	9.310	1068	1075	1079	rVV4	7401	14702	0.41%	0.034%
31	9.369	1079	1085	1090	rVV2	52384	93442	2.58%	0.215%
32	9.416	1090	1093	1097	rVV5	10696	18084	0.50%	0.042%
33	9.475	1097	1103	1112	rVB2	46983	88624	2.45%	0.204%
34	9.599	1118	1124	1130	rBV4	19726	32533	0.90%	0.075%
35	9.669	1130	1136	1141	rVB10	7256	16342	0.45%	0.038%
36	9.740	1141	1148	1161	rVB	299077	533893	14.75%	1.231%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

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_P\Methods\SFAM-EPA-BP101122.M
Title : SVOA CALIBRATION

37	9.869	1161	1170	1178	rBV10	30146	102503	2.83%	0.236%
38	10.052	1193	1201	1212	rBV10	13415	36845	1.02%	0.085%
39	10.146	1212	1217	1218	rBV3	14328	19178	0.53%	0.044%
40	10.181	1218	1223	1233	rVB3	41921	112266	3.10%	0.259%
41	10.275	1233	1239	1260	rBV	497144	934236	25.81%	2.154%
42	10.434	1260	1266	1277	rBV	127149	265914	7.35%	0.613%
43	10.599	1289	1294	1297	rBV	51485	85797	2.37%	0.198%
44	10.652	1297	1303	1318	rVB	1204354	2081813	57.52%	4.799%
45	10.799	1318	1328	1342	rBV	351208	671389	18.55%	1.548%
46	10.910	1343	1347	1351	rVV4	19076	31435	0.87%	0.072%
47	10.963	1352	1356	1362	rVV5	18058	34910	0.96%	0.080%
48	11.016	1362	1365	1371	rVB7	10821	18754	0.52%	0.043%
49	11.069	1371	1374	1378	rBV6	8302	13447	0.37%	0.031%
50	11.469	1433	1442	1444	rBV3	19273	52357	1.45%	0.121%
51	11.640	1463	1471	1476	rBV10	12695	29672	0.82%	0.068%
52	11.863	1505	1509	1514	rVB4	19267	31612	0.87%	0.073%
53	11.999	1524	1532	1538	rBV	125914	225319	6.23%	0.519%
54	12.163	1555	1560	1565	rBV3	59102	112166	3.10%	0.259%
55	12.210	1566	1568	1574	rVV3	25273	56355	1.56%	0.130%
56	12.275	1575	1579	1588	rVV8	32088	95353	2.63%	0.220%
57	12.375	1592	1596	1601	rVV3	16258	30507	0.84%	0.070%
58	12.422	1601	1604	1611	rVV8	10716	23364	0.65%	0.054%
59	12.540	1621	1624	1629	rVB3	12946	18979	0.52%	0.044%
60	12.657	1639	1644	1650	rBV2	35399	60289	1.67%	0.139%
61	12.751	1656	1660	1670	rVB10	21494	51392	1.42%	0.118%
62	13.222	1738	1740	1747	rBV8	10700	15422	0.43%	0.036%
63	13.298	1749	1753	1757	rVV5	19601	34566	0.96%	0.080%
64	13.340	1758	1760	1763	rVV3	15344	20795	0.57%	0.048%
65	13.387	1764	1768	1774	rVB	66651	105797	2.92%	0.244%
66	13.451	1777	1779	1785	rVB7	11703	16804	0.46%	0.039%
67	13.551	1791	1796	1799	rBV4	31095	48322	1.34%	0.111%
68	13.840	1839	1845	1847	rBV7	19480	38633	1.07%	0.089%
69	13.910	1851	1857	1866	rVB	856223	1220748	33.73%	2.814%
70	14.051	1877	1881	1885	rBV5	27377	46852	1.29%	0.108%
71	14.187	1898	1904	1912	rBV	984287	1520991	42.02%	3.506%
72	14.304	1920	1924	1936	rVB	47709	94147	2.60%	0.217%
73	14.422	1939	1944	1949	rVV	137865	190007	5.25%	0.438%
74	14.498	1950	1957	1963	rVV	1595768	2414708	66.71%	5.567%
75	14.557	1963	1967	1974	rVB	307230	446798	12.34%	1.030%
76	14.722	1991	1995	2001	rBV2	37547	74141	2.05%	0.171%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

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_P\Methods\SFAM-EPA-BP101122.M
Title : SVOA CALIBRATION

77	15.245	2079	2084	2090	rBV	225001	346920	9.58%	0.800%
78	15.492	2120	2126	2131	rBV	1380713	1963375	54.25%	4.526%
79	15.616	2142	2147	2152	rBV2	191022	285398	7.89%	0.658%
80	15.722	2162	2165	2166	rBV	24581	27395	0.76%	0.063%
81	15.751	2167	2170	2179	rVB3	90595	152608	4.22%	0.352%
82	16.016	2208	2215	2227	rBV	835610	1199134	33.13%	2.764%
83	16.192	2237	2245	2253	rBV2	87795	206059	5.69%	0.475%
84	16.345	2268	2271	2274	rBV	35793	39712	1.10%	0.092%
85	16.528	2297	2302	2307	rBV	461821	619052	17.10%	1.427%
86	16.798	2344	2348	2358	rVB	73364	130308	3.60%	0.300%
87	17.063	2391	2393	2399	rVB2	50238	67459	1.86%	0.156%
88	17.181	2411	2413	2418	rVB5	32255	34940	0.97%	0.081%
89	17.257	2420	2426	2435	rVV	1802483	2637690	72.88%	6.081%
90	17.351	2437	2442	2452	rVB	1429862	2079218	57.45%	4.793%
91	17.922	2537	2539	2544	rVB	42478	44331	1.22%	0.102%
92	18.139	2573	2576	2583	rBV	124473	170887	4.72%	0.394%
93	18.898	2702	2705	2709	rVB2	75043	91937	2.54%	0.212%
94	19.051	2729	2731	2736	rVB3	68256	74124	2.05%	0.171%
95	19.375	2783	2786	2795	rBV3	108675	152625	4.22%	0.352%
96	19.592	2818	2823	2827	rBV	1948110	2601312	71.87%	5.997%
97	19.639	2827	2831	2840	rVB	1373250	1660505	45.88%	3.828%
98	21.345	3116	3121	3127	rBV	2712956	3349079	92.53%	7.721%
99	23.604	3499	3505	3519	rVB2	1485426	3040689	84.01%	7.010%
100	23.763	3526	3532	3544	rVB2	1822003	3619444	100.00%	8.344%

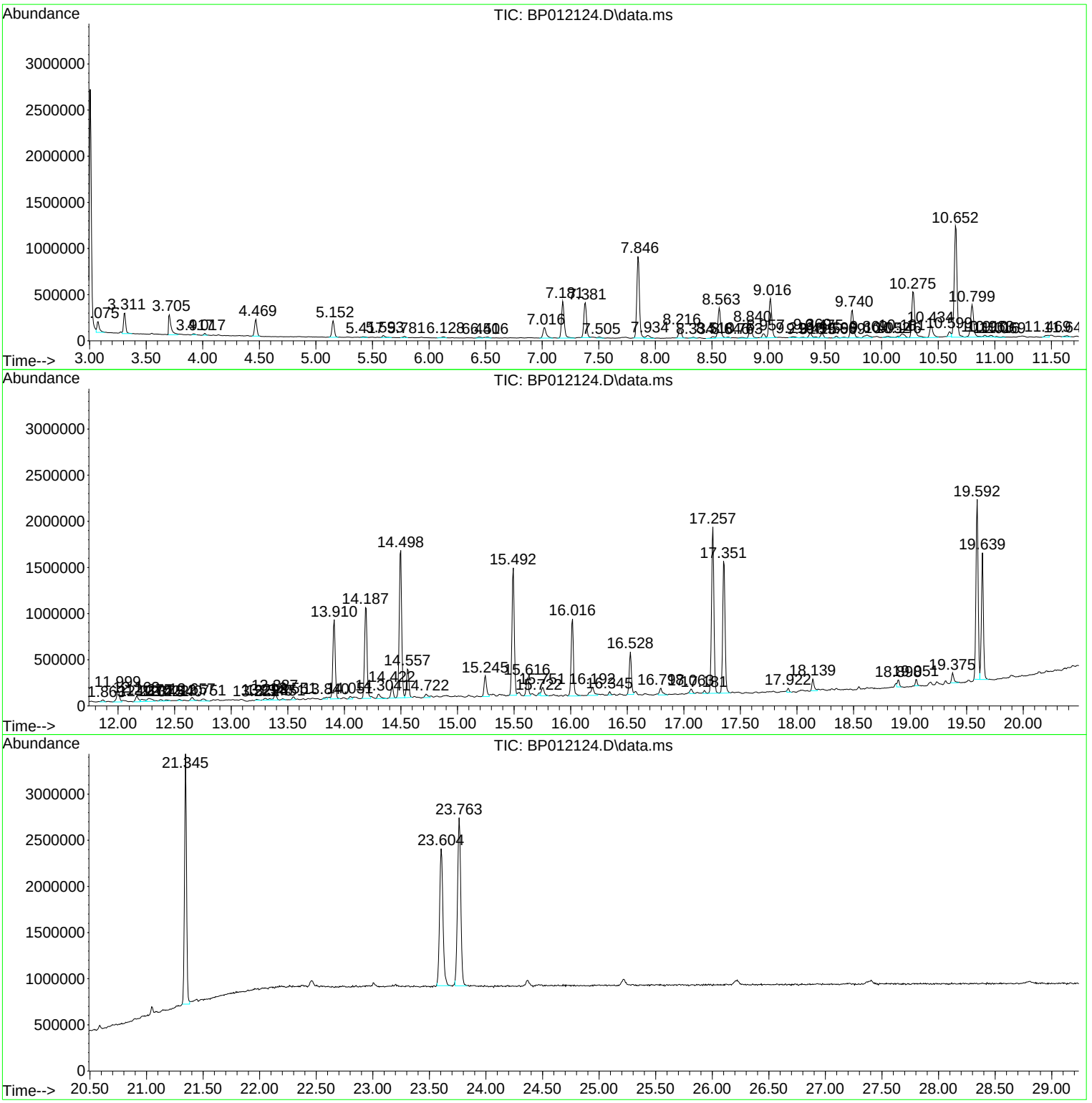
Sum of corrected areas: 43376948

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

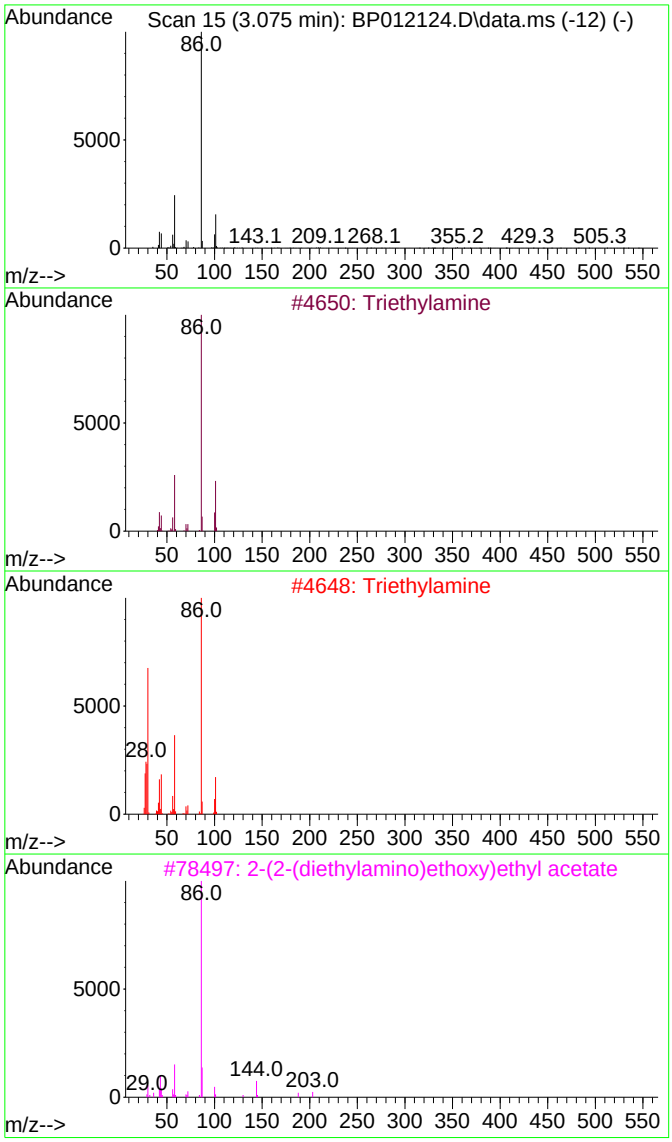
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

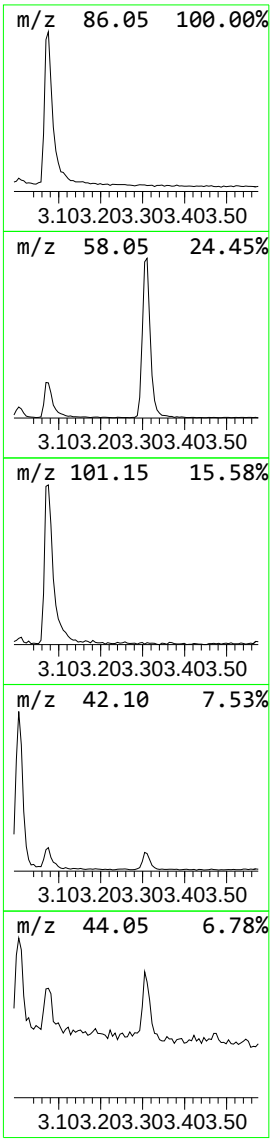
****

Peak Number 1 Triethylamine Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	2.40 ng/ul	171727	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylamine	101	C6H15N	000121-44-8	90
2	Triethylamine	101	C6H15N	000121-44-8	83
3	2-(2-(diethylamino)ethoxy)ethyl ...	203	C10H21NO3	1000456-44-8	72
4	Triethylamine	101	C6H15N	000121-44-8	52
5	Triethylamine	101	C6H15N	000121-44-8	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

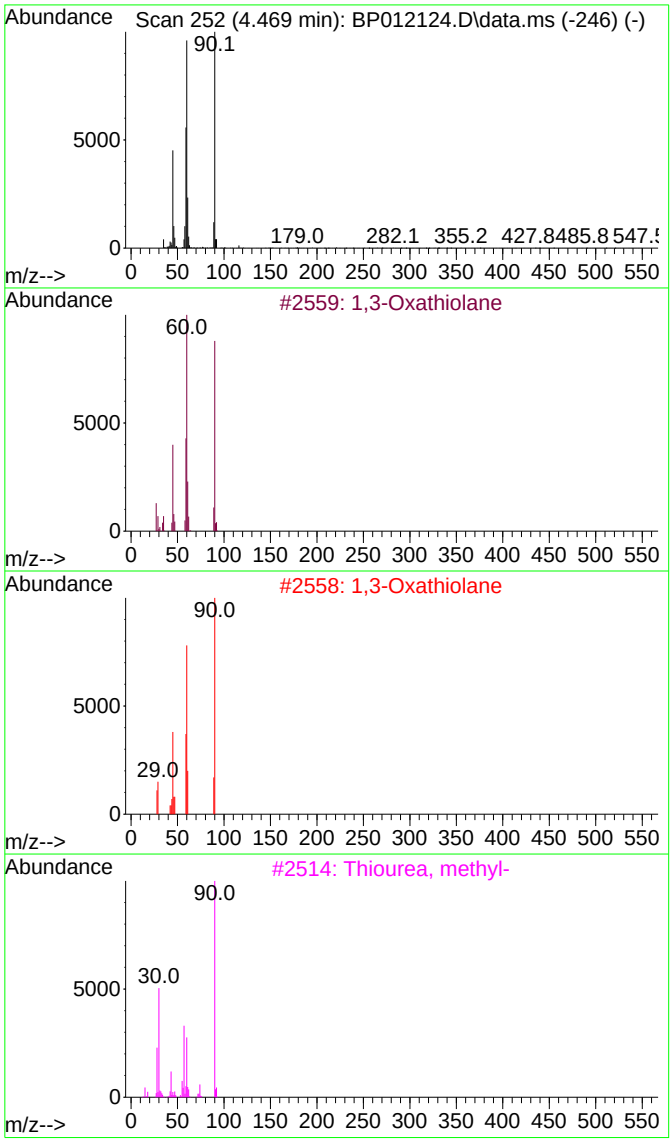
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

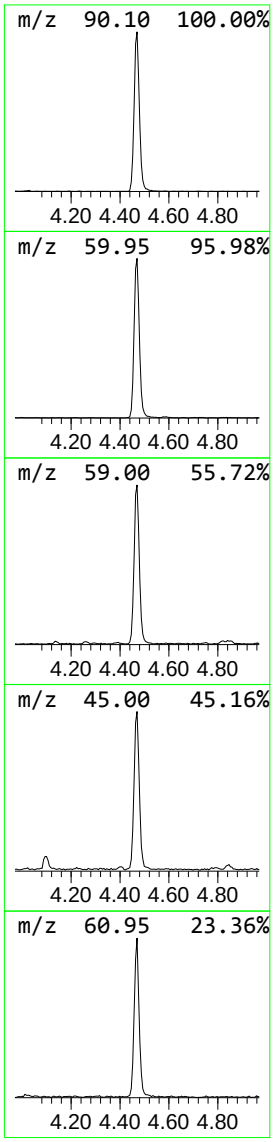
****

Peak Number 2 1,3-Oxathiolane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.469	3.69 ng/ul	263905	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	94
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
3	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
4	3-Thietanol	90	C3H6OS	010304-16-2	42
5	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

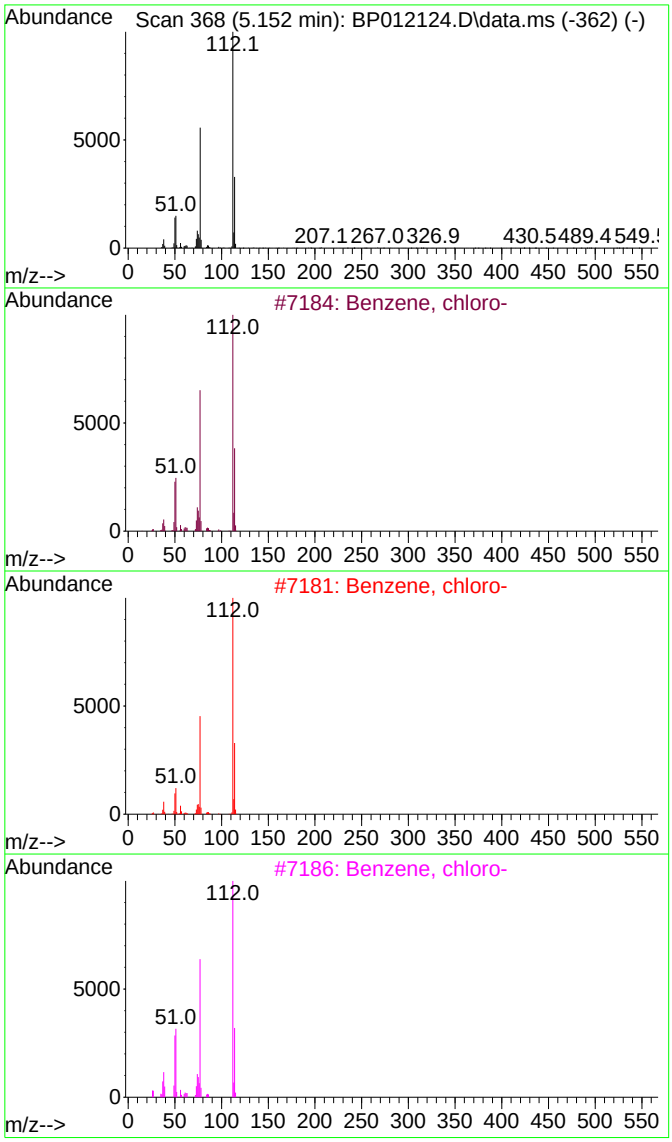
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

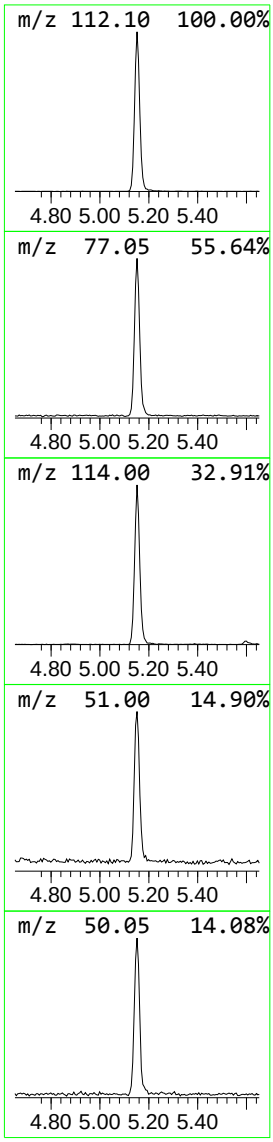
****

Peak Number 3 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.152	3.69 ng/ul	263662	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

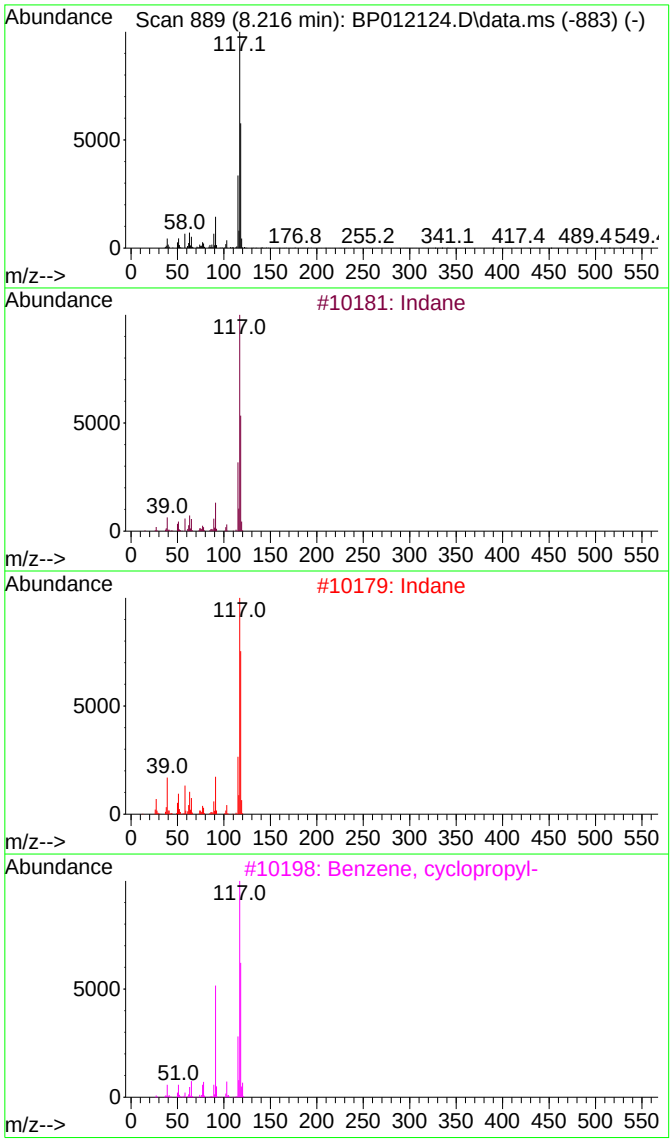
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

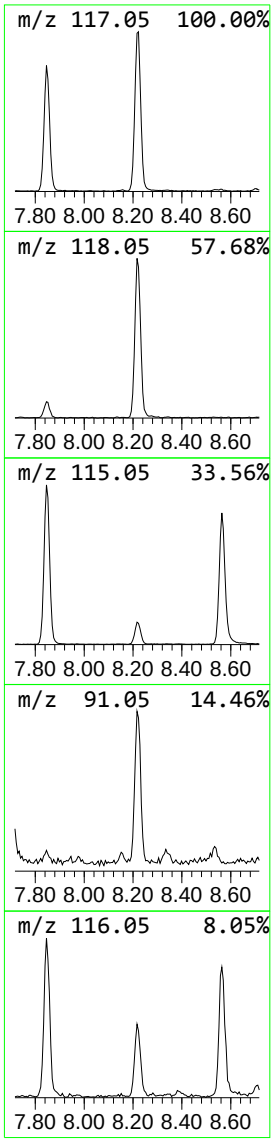
****

Peak Number 4 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.216	2.62 ng/ul	187499	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Indane	118	C9H10	000496-11-7	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

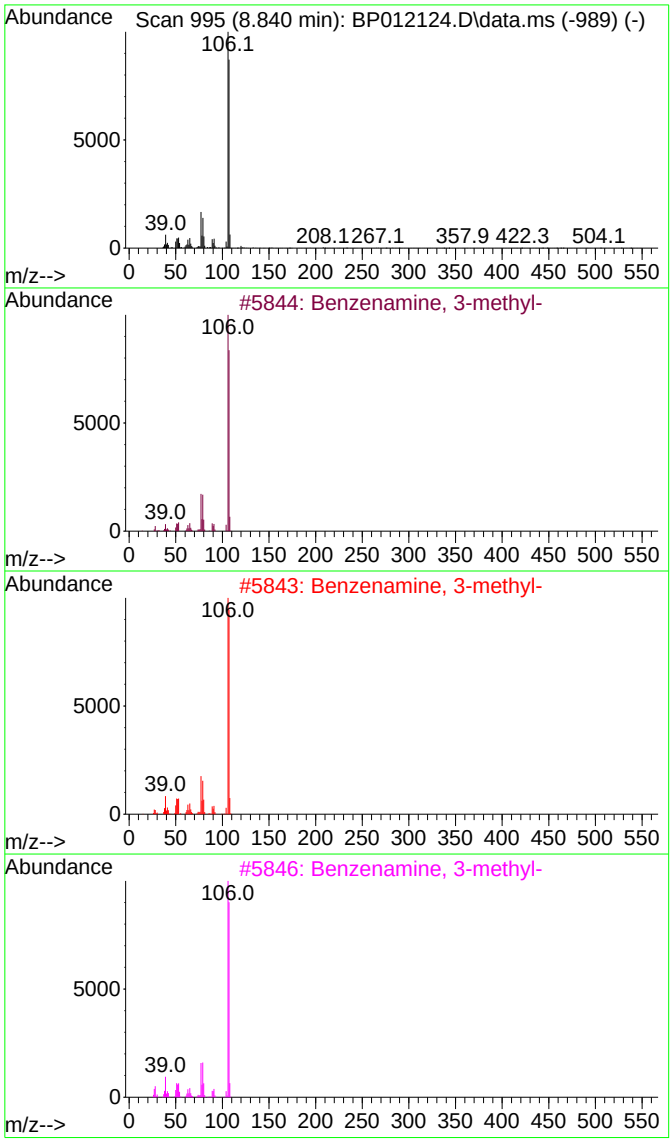
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

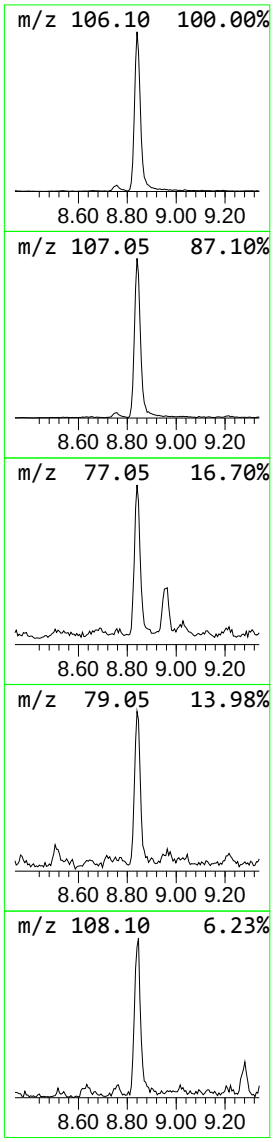
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzenamine, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.840	3.48 ng/ul	248702	1,4-Dichlorobenzene-d4	7.846

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94
5	o-Toluidine	107	C7H9N	000095-53-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

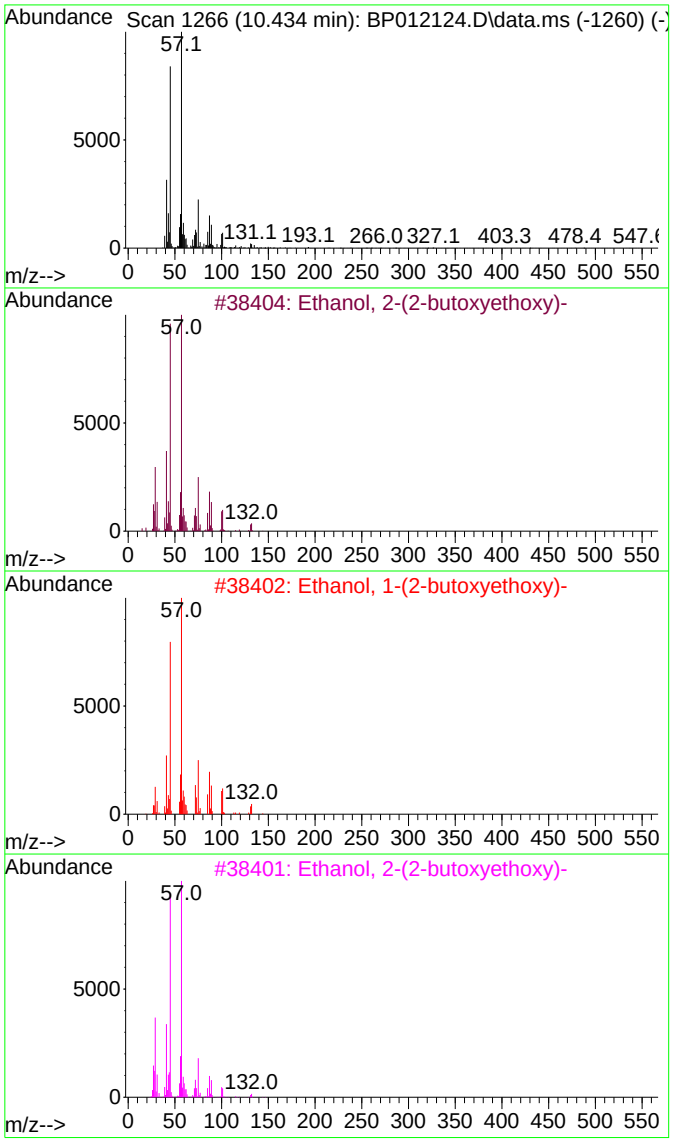
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

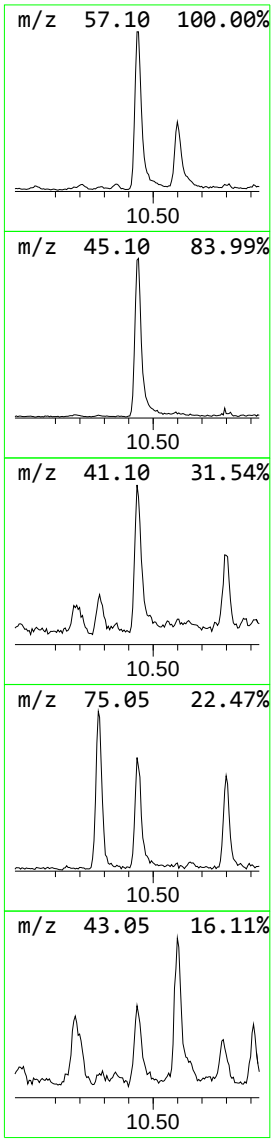
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.434	2.55 ng/ul	265914	Naphthalene-d8	10.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

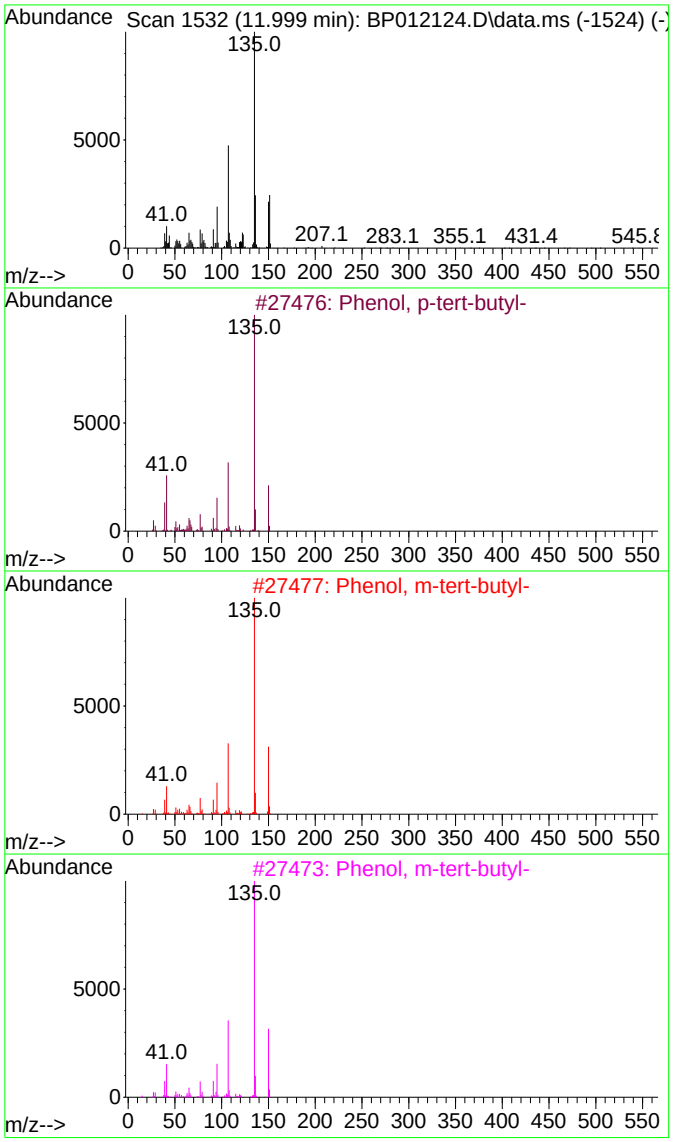
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

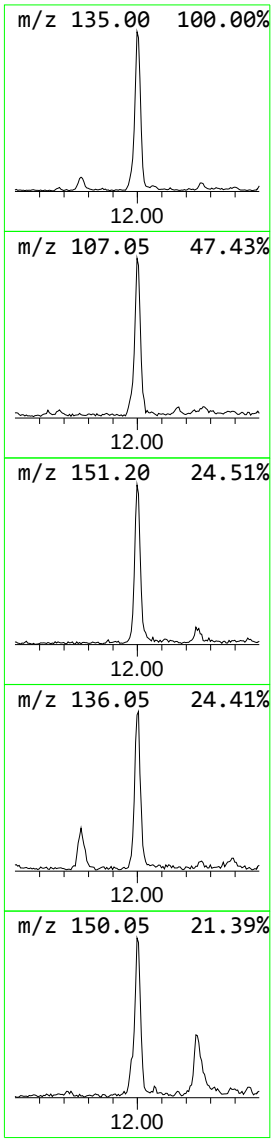
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Phenol, p-tert-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.999	2.16 ng/ul	225319	Naphthalene-d8	10.652

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	59
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	58
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	58





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Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

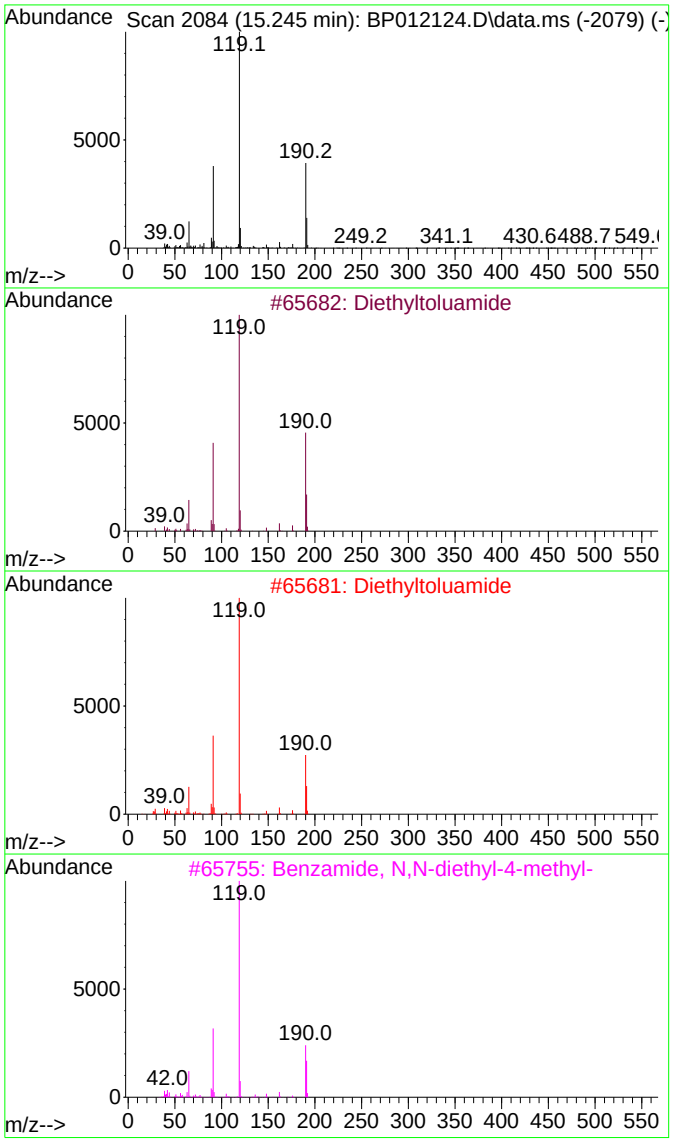
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

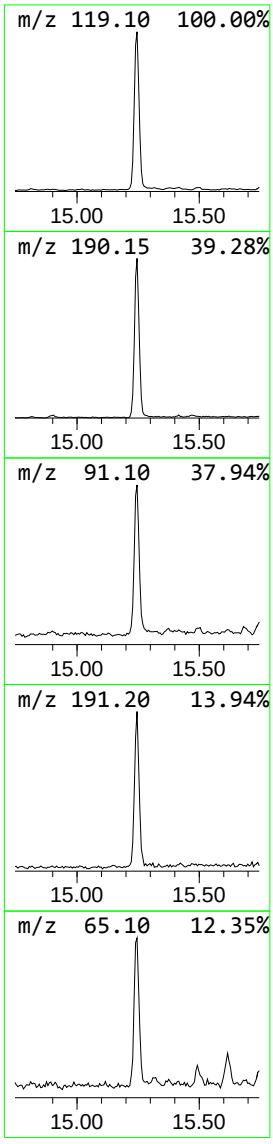
****

Peak Number 8 Diethyltoluamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.245	2.87 ng/ul	346920	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	95
3	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	76
4	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76
5	Benzamide, 4-methyl-N-butyl-	191	C12H17NO	1000407-46-6	68





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

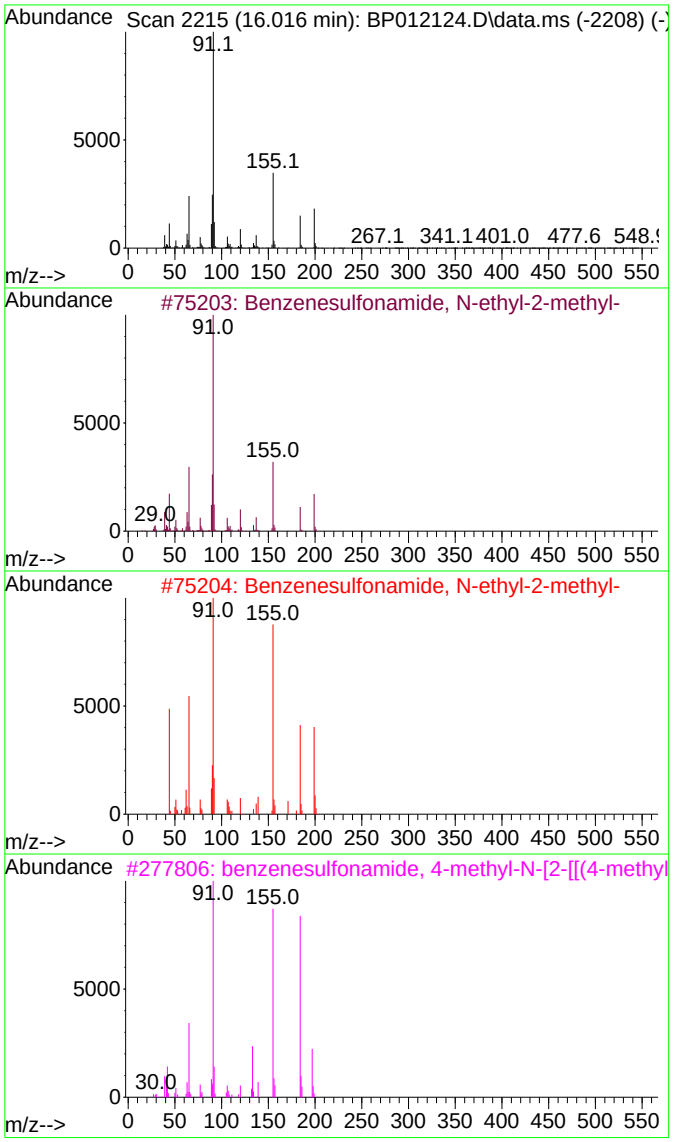
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

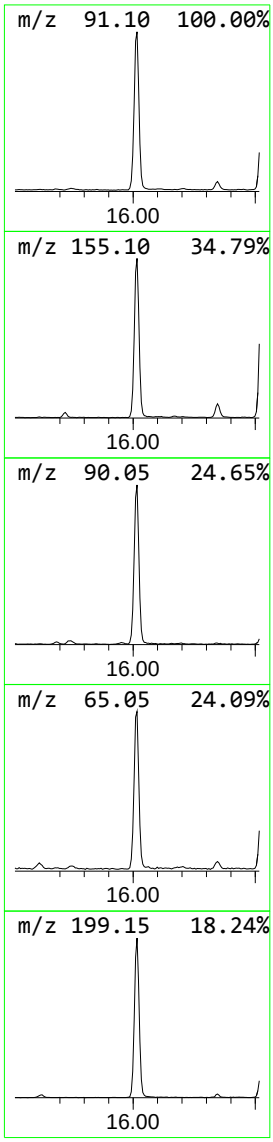
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzenesulfonamide, N-ethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.016	9.09 ng/ul	1199130	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	97
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	53
3	benzenesulfonamide, 4-methyl-N-[...	369	C16H19NO5S2	1000402-61-7	50
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	Benzenesulfonyl chloride, 4-methyl-	190	C7H7ClO2S	000098-59-9	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

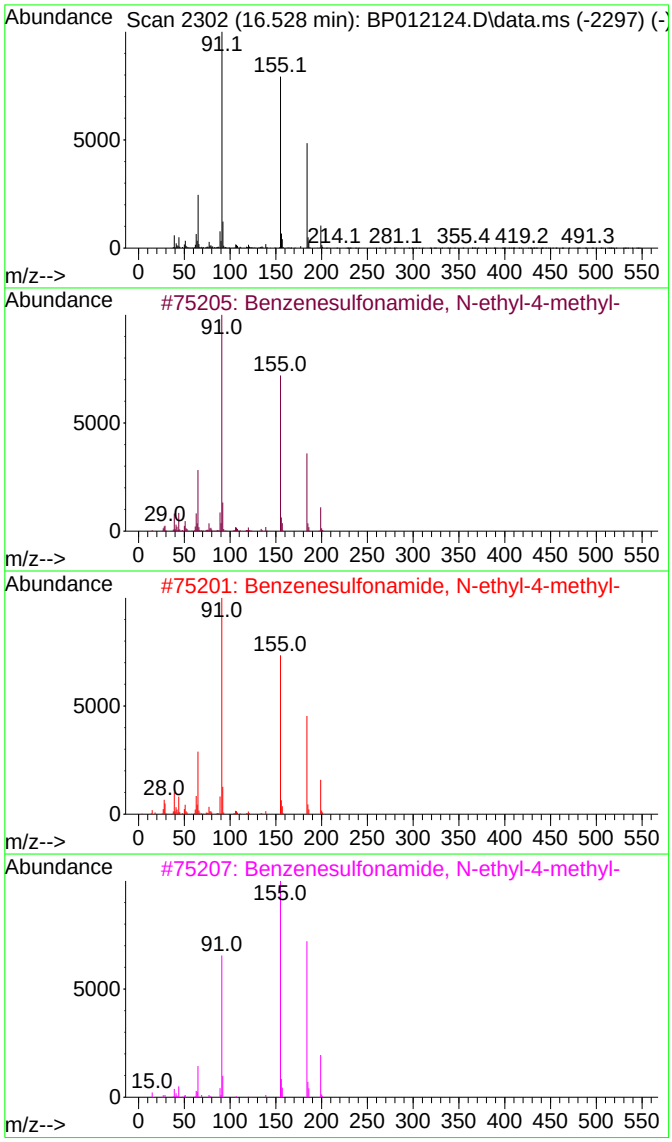
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

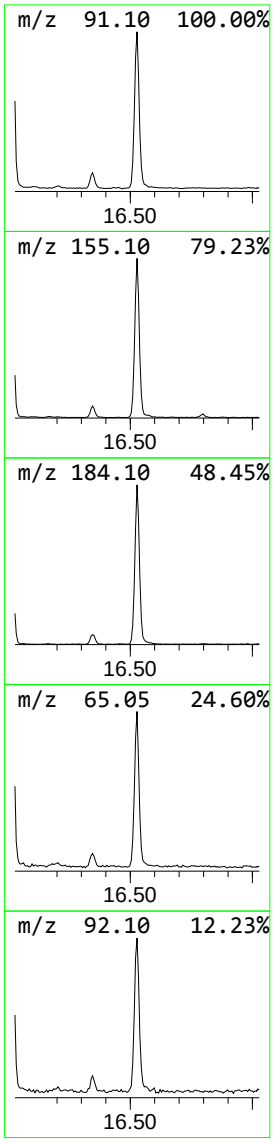
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.528	4.69 ng/ul	619052	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	94
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	87
5	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012124.D
Acq On : 13 Oct 2022 17:04
Operator : CG/JU
Sample : N4976-12DL 2X
Misc :
ALS Vial : 95 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.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
Triethylamine	3.075	2.4	ng/ul	171727	1	7.846	1430260	20.0
1,3-Oxathiolane	4.469	3.7	ng/ul	263905	1	7.846	1430260	20.0
Benzene, chloro-	5.152	3.7	ng/ul	263662	1	7.846	1430260	20.0
Indane	8.216	2.6	ng/ul	187499	1	7.846	1430260	20.0
Benzenamine, 3-...	8.840	3.5	ng/ul	248702	1	7.846	1430260	20.0
Ethanol, 2-(2-b...	10.434	2.5	ng/ul	265914	2	10.652	2081810	20.0
Phenol, p-tert-...	11.999	2.2	ng/ul	225319	2	10.652	2081810	20.0
Diethyltoluamide	15.245	2.9	ng/ul	346920	3	14.498	2414710	20.0
Benzenesulfonam...	16.016	9.1	ng/ul	1199130	4	17.257	2637690	20.0
Benzenesulfonam...	16.528	4.7	ng/ul	619052	4	17.257	2637690	20.0