LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BHA48

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_N\Methods\SFAM-EPA-BN101222.M
Title : SVOA CALIBRATION

Signal : TIC: BN022372.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	5	8	23	rVB	155891	282349	7.52%	0.543%
2	3.246	40	44	48	rBV	44975	58070	1.55%	0.112%
3	3.699	118	121	135	rBV	55033	99331	2.65%	0.191%
4	3.922	153	159	166	rBV	32904	56839	1.51%	0.109%
5	5.187	367	374	391	rBV	2078764	3233710	86.11%	6.224%
6	5.322	392	397	403	rVV	18812	29895	0.80%	0.058%
7	5.605	438	445	450	rBV2	26321	40567	1.08%	0.078%
8	6.393	570	579	585	rBV	80334	138262	3.68%	0.266%
9	6.634	616	620	627	rVV	20735	31712	0.84%	0.061%
10	6.893	656	664	674	rBV3	22669	48439	1.29%	0.093%
11	7.093	691	698	710	rVV	207376	349841	9.32%	0.673%
12	7.263	717	727	734	rBV	773674	1229444	32.74%	2.366%
13	7.357	738	743	753	rVB	44318	69612	1.85%	0.134%
14	7.457	753	760	768	rBV	732281	1199795	31.95%	2.309%
15	7.522	768	771	778	rVV	40544	66461	1.77%	0.128%
16	7.628	783	789	795	rVB	51516	76898	2.05%	0.148%
17	7.934	830	841	854	rBV	658645	1121482	29.86%	2.159%
18	8.128	870	874	880	rVV	55754	85664	2.28%	0.165%
19	8.304	896	904	914	rVB2	38276	82811	2.21%	0.159%
20	8.652	956	963	973	rBV	531693	875927	23.33%	1.686%
21	8.752	974	980	990	rVB8	10994	31192	0.83%	0.060%
22	8.852	990	997	1005	rVB8	16673	30811	0.82%	0.059%
23	8.940	1005	1012	1020	rBV5	28459	56006	1.49%	0.108%
24	9.116	1035	1042	1049	rBV	846483	1400048	37.28%	2.695%
25	9.840	1159	1165	1178	rVB	654276	1093783	29.13%	2.105%
26	9.999	1186	1192	1196	rBV3	16399	28634	0.76%	0.055%
27	10.181	1218	1223	1231	rVV2	21012	45718	1.22%	0.088%
28	10.381	1250	1257	1267	rBV	906976	1566514	41.72%	3.015%
29	10.463	1267	1271	1273	rVV3	21616	36931	0.98%	0.071%
30	10.498	1274	1277	1279	rVV4	28828	45093	1.20%	0.087%
31	10.534	1279	1283	1303	rVB	212203	481455	12.82%	0.927%
32	10.763	1314	1322	1329	rBV	822074	1392414	37.08%	2.680%
33	10.851	1329	1337	1340	rVV	117262	210734	5.61%	0.406%
34	10.898	1340	1345	1361	rVV2	410453	796972	21.22%	1.534%
35	11.134	1377	1385	1396	rVB9	9545	28248	0.75%	0.054%
36	11.363	1418	1424	1432	rVB3	20042	39244	1.05%	0.076%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : BHA48

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_N\Methods\SFAM-EPA-BN101222.M
Title : SVOA CALIBRATION

37	11.916	1506	1518	1523	rBV2	36932	107003	2.85%	0.206%
38	11.963	1523	1526	1534	rVB7	16468	36142	0.96%	0.070%
39	12.110	1541	1551	1558	rBV	232054	462771	12.32%	0.891%
40	12.181	1559	1563	1573	rVV2	21094	43186	1.15%	0.083%
41	12.293	1575	1582	1586	rVV7	13091	30373	0.81%	0.058%
42	12.357	1589	1593	1597	rVV	21115	40467	1.08%	0.078%
43	12.581	1623	1631	1641	rBV10	16083	48980	1.30%	0.094%
44	13.140	1721	1726	1731	rVV	91489	128208	3.41%	0.247%
45	13.204	1731	1737	1740	rVV6	18281	37784	1.01%	0.073%
46	13.240	1740	1743	1751	rVB	37802	56821	1.51%	0.109%
47	13.416	1766	1773	1778	rBV3	16279	29347	0.78%	0.056%
48	13.663	1809	1815	1819	rBV	51008	80425	2.14%	0.155%
49	14.022	1869	1876	1888	rVV	1521216	2188635	58.28%	4.213%
50	14.139	1893	1896	1900	rVB	25654	32065	0.85%	0.062%
51	14.304	1918	1924	1932	rVV	1794063	2671914	71.15%	5.143%
52	14.392	1934	1939	1944	rVB	56063	80036	2.13%	0.154%
53	14.610	1969	1976	1983	rBV	1153293	1723612	45.90%	3.318%
54	14.822	2006	2012	2025	rBV	128441	234356	6.24%	0.451%
55	15.186	2068	2074	2084	rVB	41268	93713	2.50%	0.180%
56	15.351	2098	2102	2106	rBV	37412	47187	1.26%	0.091%
57	15.504	2122	2128	2133	rVV5	17974	33508	0.89%	0.064%
58	15.604	2138	2145	2157	rVB	2220322	3168325	84.37%	6.098%
59	15.728	2158	2166	2174	rBV	700750	973084	25.91%	1.873%
60	15.822	2178	2182	2184	rVV	23462	29393	0.78%	0.057%
61	15.869	2184	2190	2202	rVB	2106038	2805636	74.71%	5.400%
62	16.016	2211	2215	2219	rBV2	19531	28622	0.76%	0.055%
63	16.116	2225	2232	2240	rBV3	86035	159828	4.26%	0.308%
64	16.298	2258	2263	2268	rVV	105159	161953	4.31%	0.312%
65	16.657	2316	2324	2330	rVB4	81488	165767	4.41%	0.319%
66	16.739	2335	2338	2348	rVB3	26307	58857	1.57%	0.113%
67	17.116	2397	2402	2408	rBV2	28934	56029	1.49%	0.108%
68	17.275	2425	2429	2432	rBV2	27650	33011	0.88%	0.064%
69	17.322	2432	2437	2440	rBV	172965	241408	6.43%	0.465%
70	17.363	2440	2444	2454	rVB	1251601	1757265	46.79%	3.382%
71	17.463	2454	2461	2472	rBV2	2273598	3306121	88.04%	6.364%
72	17.733	2502	2507	2511	rBV5	20393	45038	1.20%	0.087%
73	17.998	2548	2552	2554	rBV	43862	56608	1.51%	0.109%
74	18.039	2554	2559	2565	rVV2	146026	277608	7.39%	0.534%
75	18.104	2567	2570	2580	rVB	31187	54313	1.45%	0.105%
76	18.257	2592	2596	2599	rBV2	29406	37839	1.01%	0.073%

LSC Area Percent Report

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Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

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

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Title : SVOA CALIBRATION

77	18.339	2605	2610	2617	rVB2	126364	181042	4.82%	0.348%
78	18.416	2618	2623	2635	rVB	172278	237476	6.32%	0.457%
79	18.516	2636	2640	2643	rBV	136104	194167	5.17%	0.374%
80	18.569	2643	2649	2654	rVB	258455	450979	12.01%	0.868%
81	19.104	2737	2740	2746	rVB	41720	57278	1.53%	0.110%
82	19.210	2755	2758	2761	rVB	55589	56207	1.50%	0.108%
83	19.269	2766	2768	2773	rVB3	42137	55237	1.47%	0.106%
84	19.327	2774	2778	2782	rBV3	44428	64354	1.71%	0.124%
85	19.375	2782	2786	2799	rVB3	84743	164401	4.38%	0.316%
86	19.480	2800	2804	2808	rBV4	41830	53279	1.42%	0.103%
87	19.769	2847	2853	2857	rBV	2605111	3500278	93.21%	6.737%
88	19.816	2857	2861	2871	rVB	282217	386252	10.29%	0.743%
89	19.986	2886	2890	2898	rVB2	59944	107832	2.87%	0.208%
90	20.063	2899	2903	2907	rBV	56213	73545	1.96%	0.142%
91	20.174	2919	2922	2926	rBV	65336	79681	2.12%	0.153%
92	20.222	2926	2930	2934	rVB	112295	134109	3.57%	0.258%
93	20.504	2973	2978	2984	rBV2	184887	250006	6.66%	0.481%
94	20.786	3023	3026	3030	rBV3	44776	56481	1.50%	0.109%
95	21.274	3105	3109	3114	rBV3	74981	116772	3.11%	0.225%
96	21.474	3140	3143	3148	rVB3	57351	66588	1.77%	0.128%
97	21.563	3153	3158	3165	rBV	1304597	1763791	46.97%	3.395%
98	21.692	3176	3180	3184	rBV	125393	170030	4.53%	0.327%
99	23.862	3542	3549	3563	rVB2	1772127	3755251	100.00%	7.228%
100	24.027	3570	3577	3587	rVB2	901919	1894989	50.46%	3.647%

Sum of corrected areas: 51954169

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
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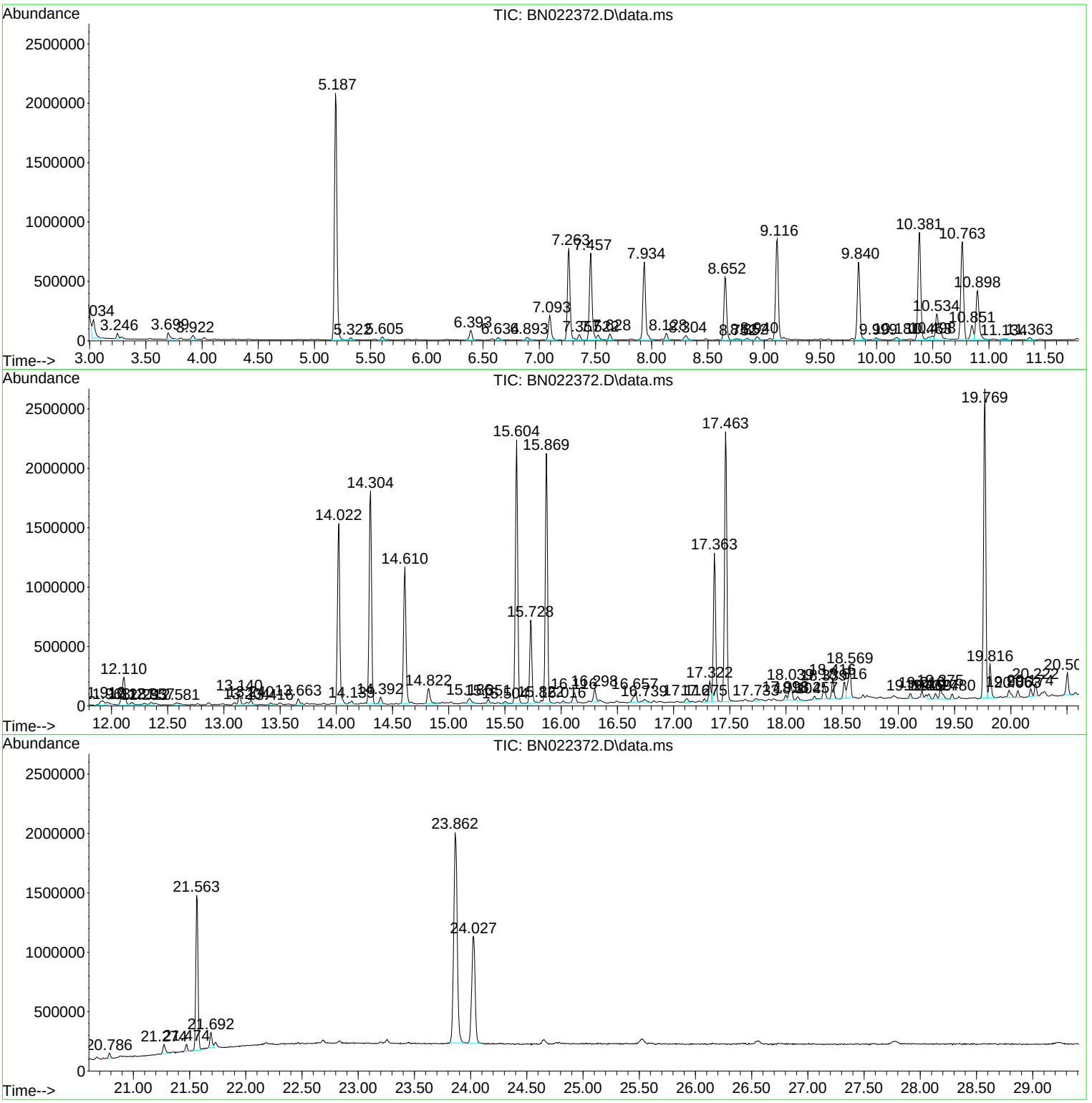
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
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Operator : CG/JU
Sample : N5232-08
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

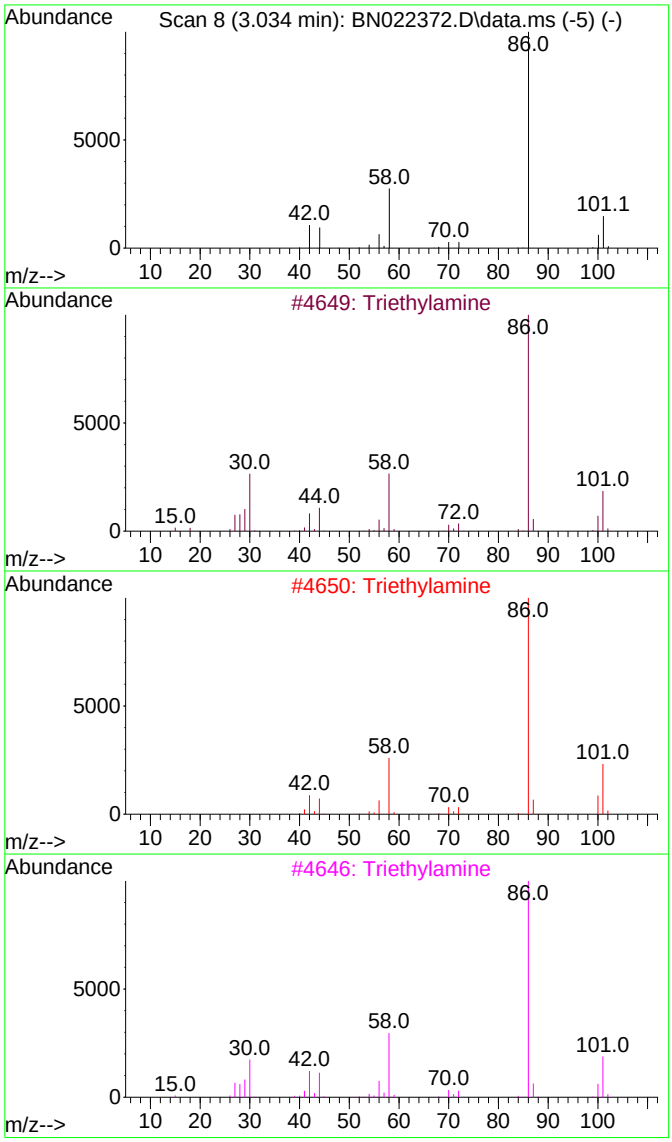
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

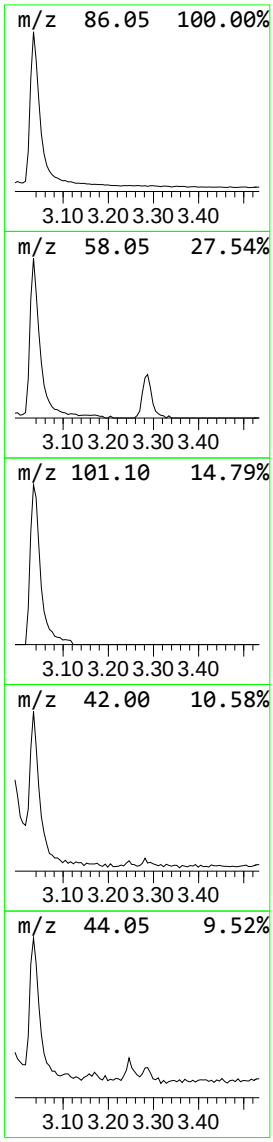
****

Peak Number 1 Triethylamine Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	5.04 ng/ul	282349	1,4-Dichlorobenzene-d4	7.934

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylamine	101	C6H15N	000121-44-8	91
2	Triethylamine	101	C6H15N	000121-44-8	90
3	Triethylamine	101	C6H15N	000121-44-8	90
4	Triethylamine	101	C6H15N	000121-44-8	86
5	Tetraethylammonium chloride	165	C8H20ClN	000056-34-8	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

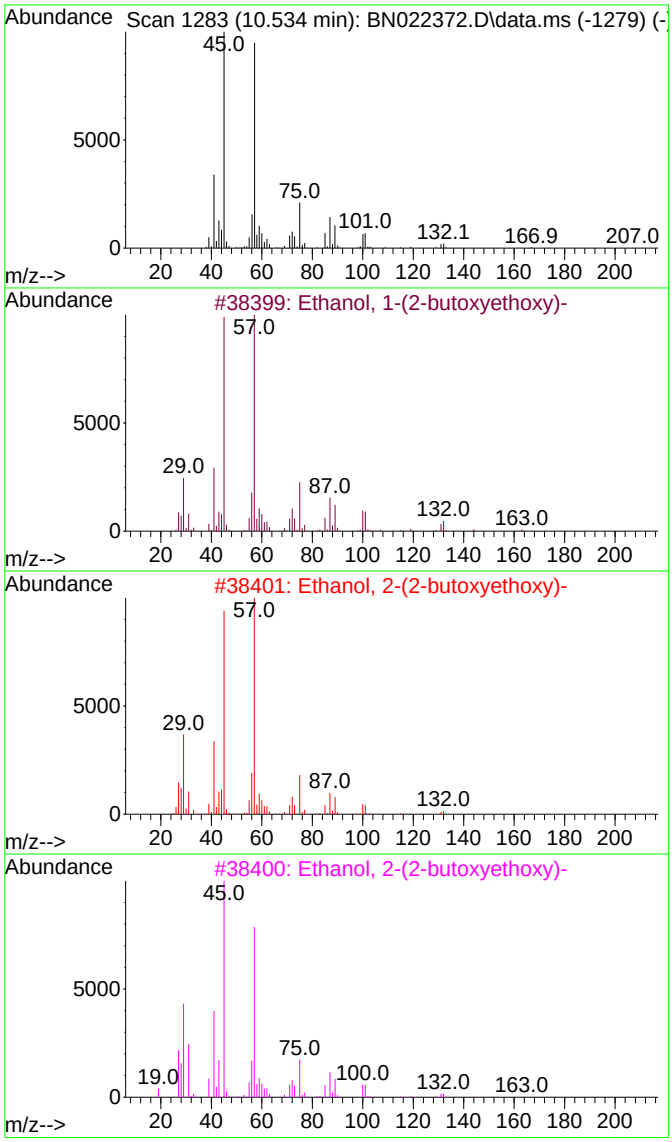
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

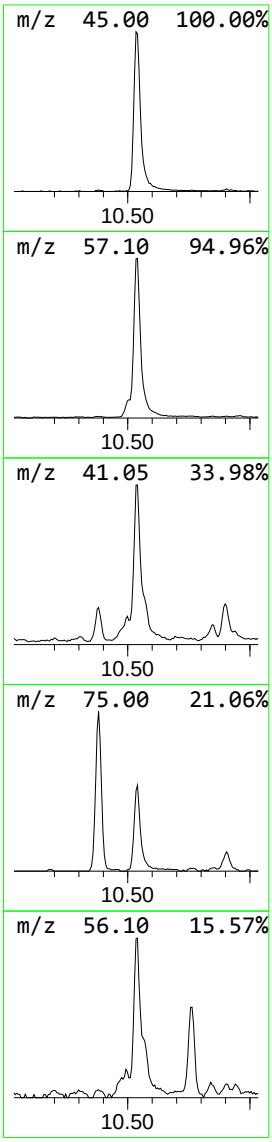
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.534	6.92 ng/ul	481455	Naphthalene-d8	10.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

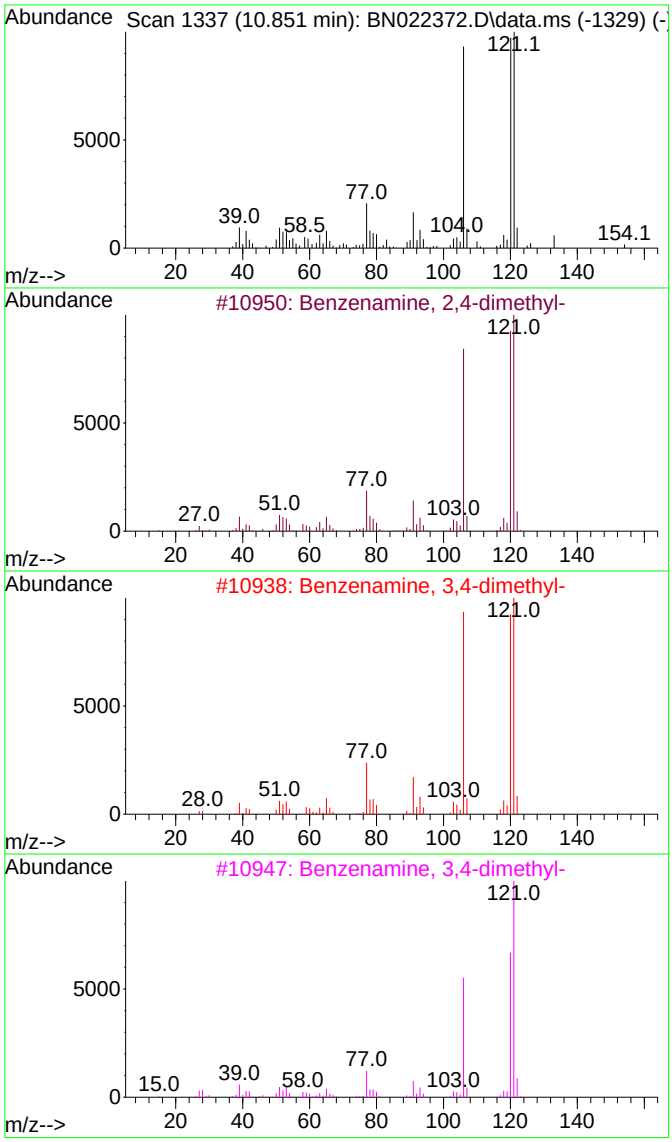
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

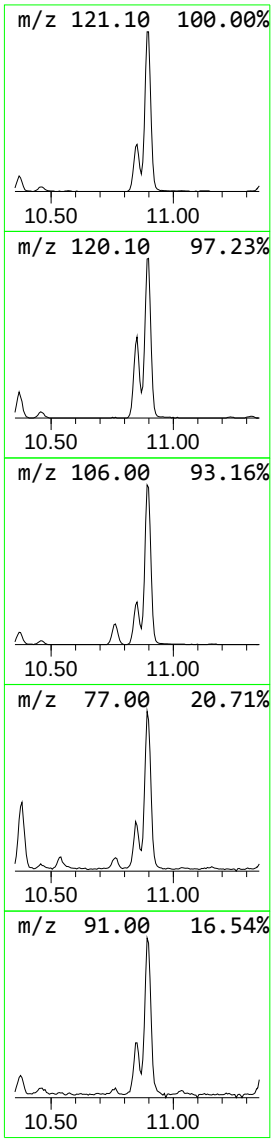
****

Peak Number 5 Benzenamine, 2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.851	3.03 ng/ul	210734	Naphthalene-d8	10.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	96
2	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
3	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
4	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
5	Benzenamine, 3,5-dimethyl-	121	C8H11N	000108-69-0	95





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Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
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Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

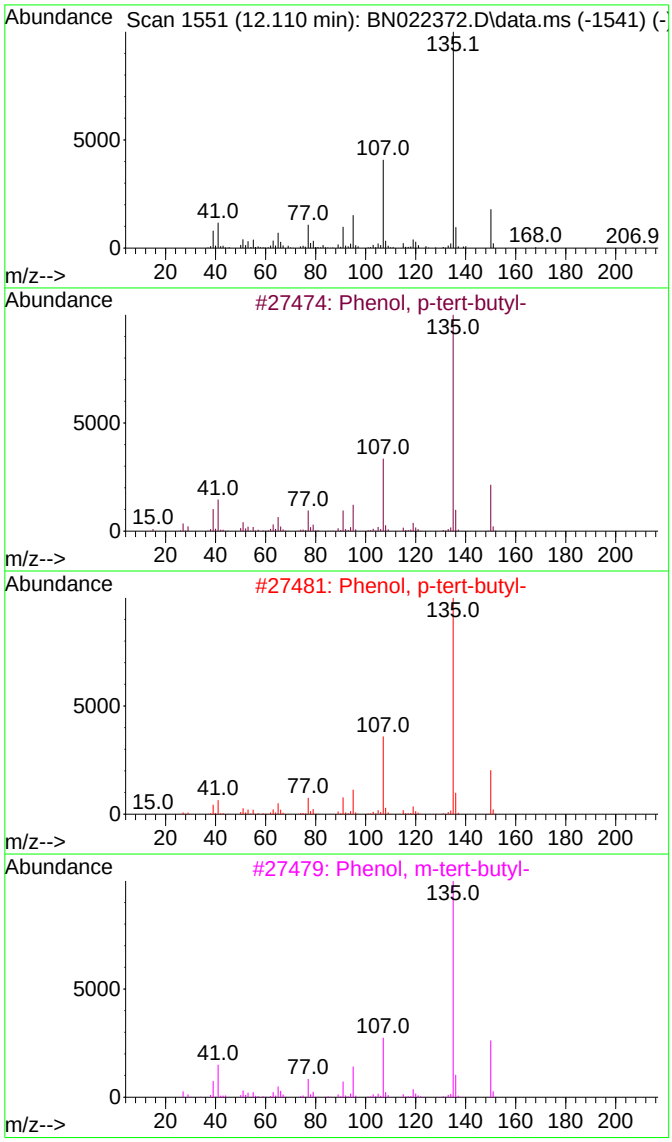
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

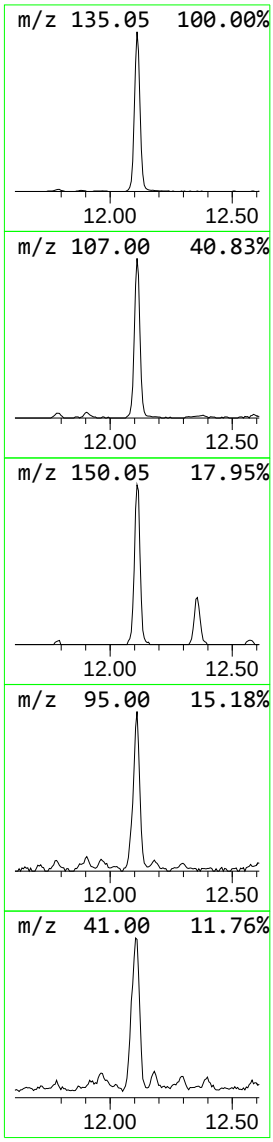
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.110	6.65 ng/ul	462771	Naphthalene-d8	10.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	95
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102922\
Data File : BN022372.D
Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

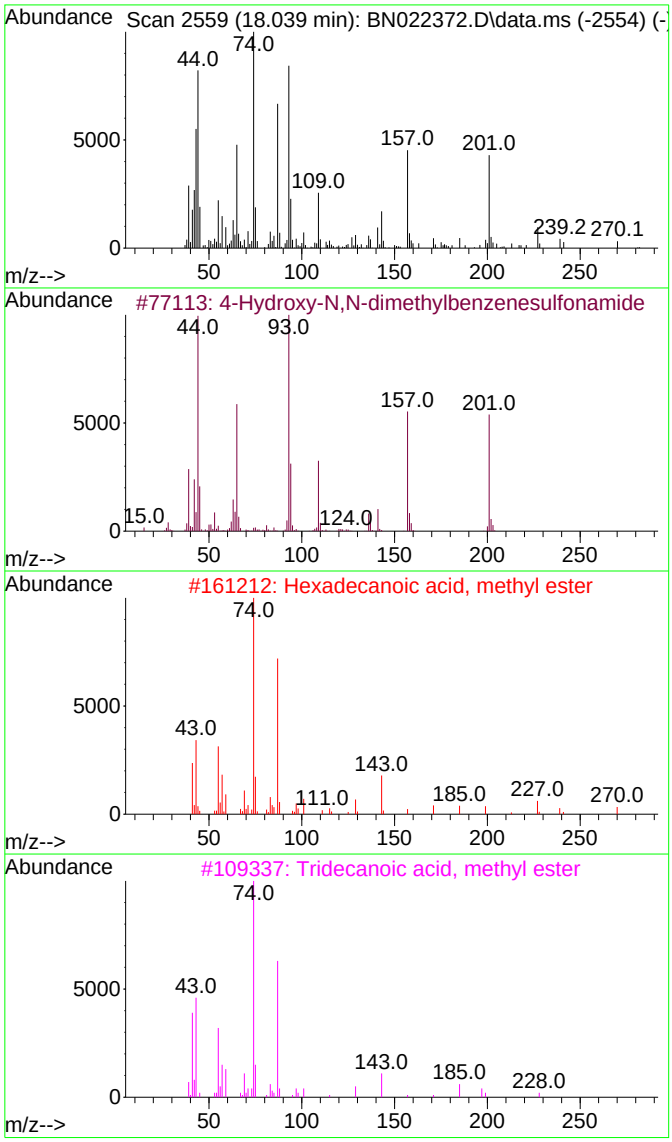
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

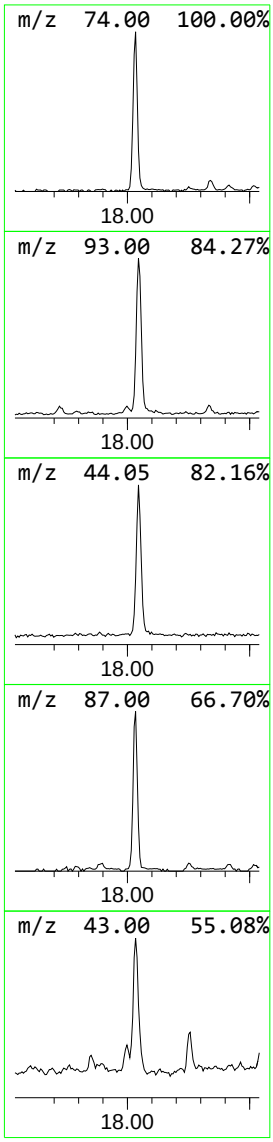
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 4-Hydroxy-N,N-dimethylbenze... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	3.16 ng/ul	277608	Phenanthrene-d10	17.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Hydroxy-N,N-dimethylbenzenesul...	201	C8H11NO3S	015020-57-2	99
2	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	56
3	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	25
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	25
5	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	15





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Acq On : 27 Oct 2022 21:48
Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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

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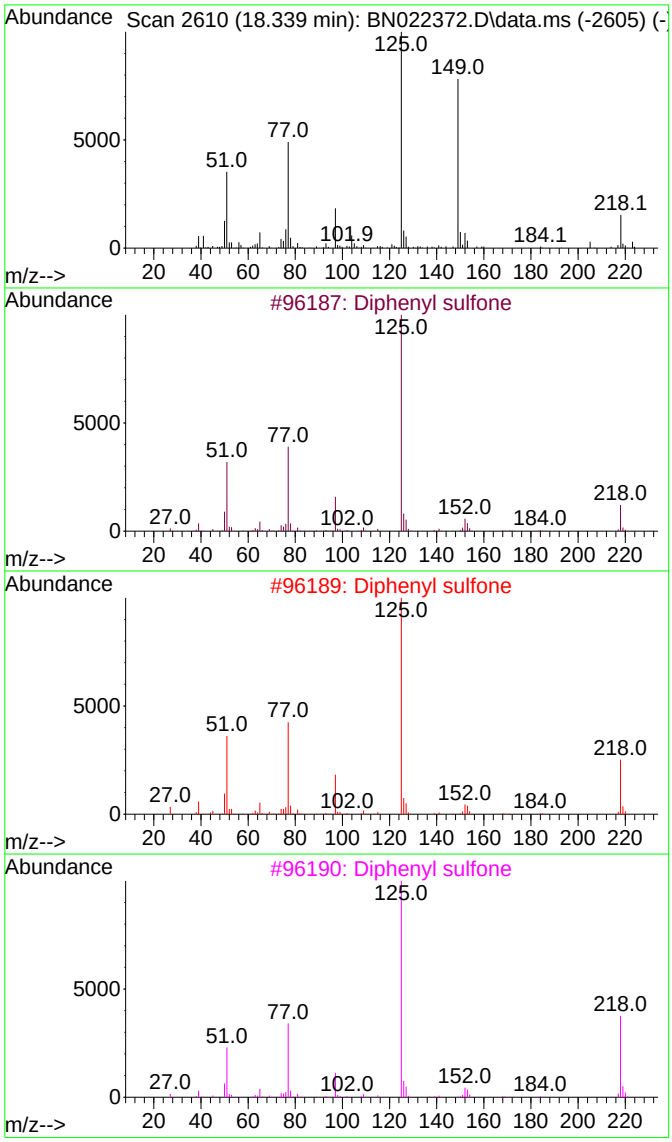
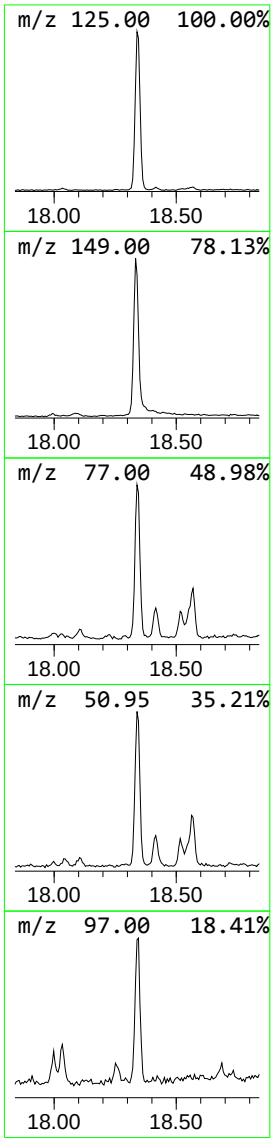
Peak Number 8 Diphenyl sulfone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.339	2.06 ng/ul	181042	Phenanthrene-d10	17.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl sulfone	218	C12H10O2S	000127-63-9	92
2	Diphenyl sulfone	218	C12H10O2S	000127-63-9	92
3	Diphenyl sulfone	218	C12H10O2S	000127-63-9	64
4	Benzene, 1-ethenyl-4-nitro-	149	C8H7NO2	000100-13-0	43
5	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	38





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Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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

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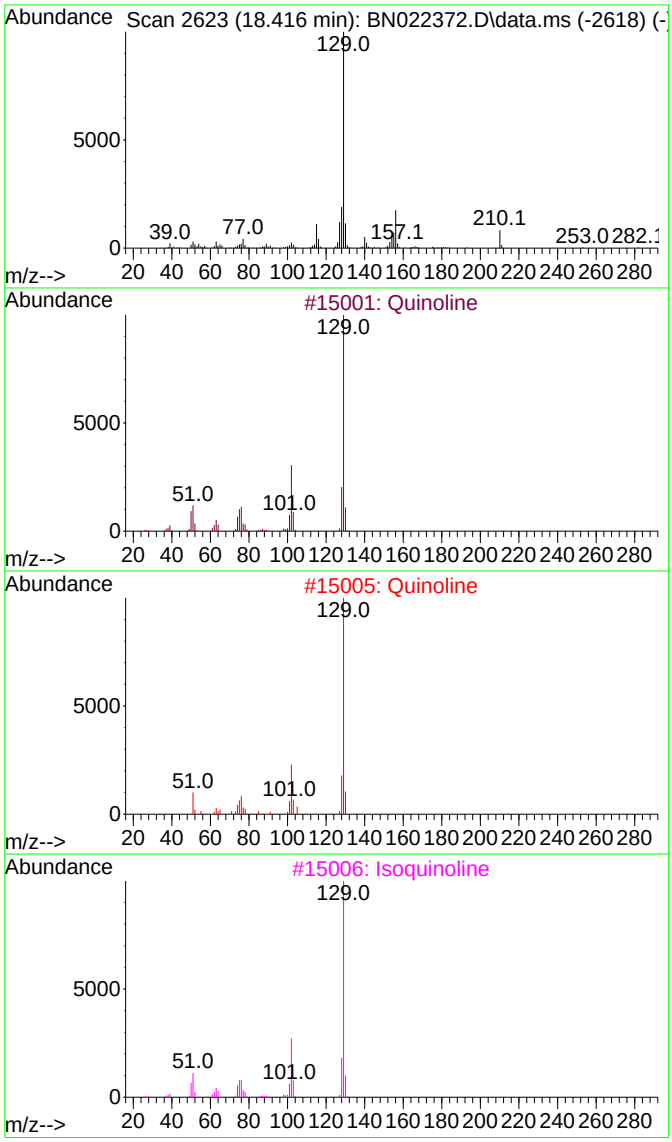
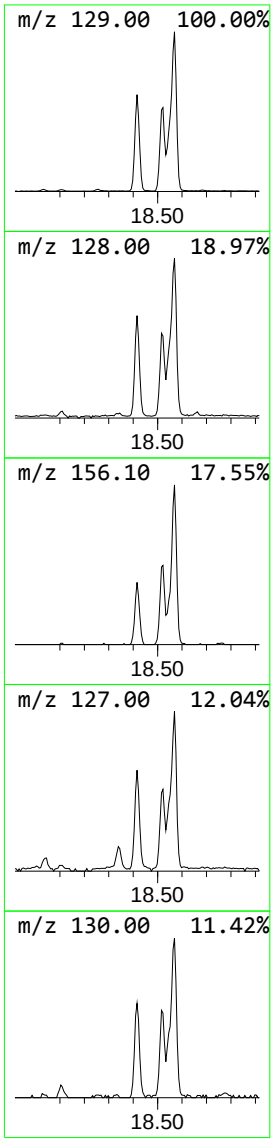
Peak Number 9 Quinoline Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.416	2.70 ng/ul	237476	Phenanthrene-d10	17.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline	129	C9H7N	000091-22-5	58
2	Quinoline	129	C9H7N	000091-22-5	53
3	Isoquinoline	129	C9H7N	000119-65-3	50
4	2-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-39-3	46
5	Naphthalene, 1,2-dihydro-2-methyl-	144	C11H12	021564-79-4	42





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Operator : CG/JU
Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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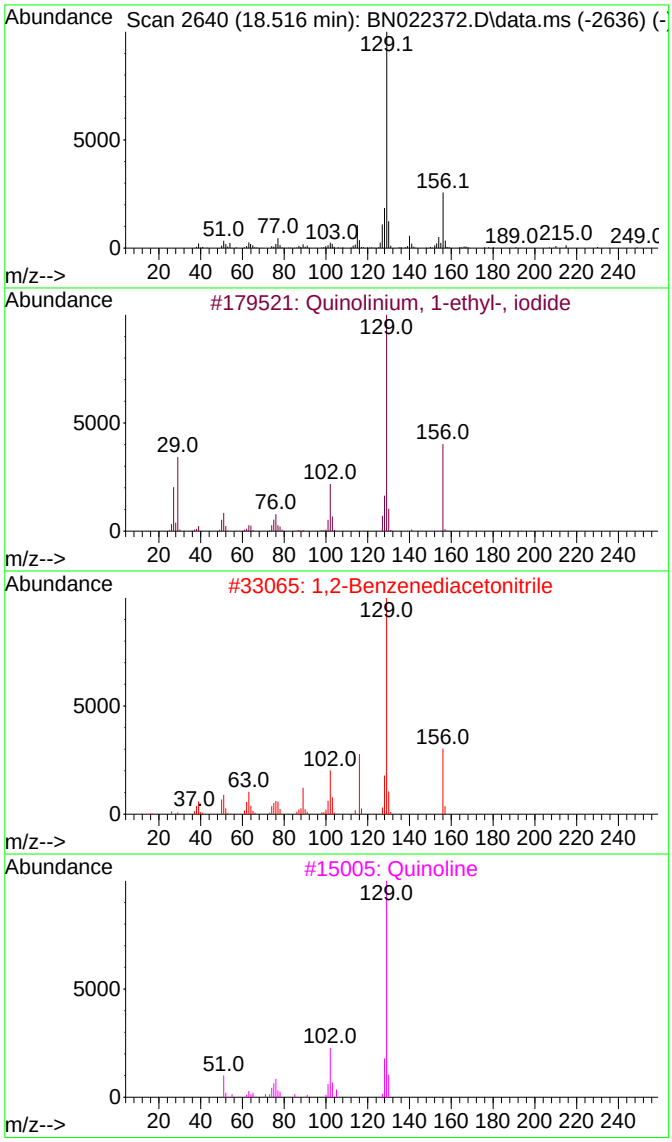
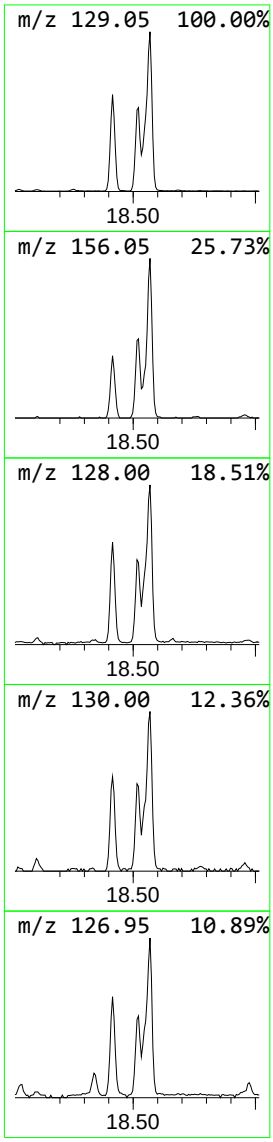
Peak Number 10 Quinolinium, 1-ethyl-, iodide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.516	2.21 ng/ul	194167	Phenanthrene-d10	17.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinolinium, 1-ethyl-, iodide	285	C11H12IN	000634-35-5	64
2	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	59
3	Quinoline	129	C9H7N	000091-22-5	50
4	Isoquinoline	129	C9H7N	000119-65-3	49
5	Isoquinoline	129	C9H7N	000119-65-3	47





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Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHA48

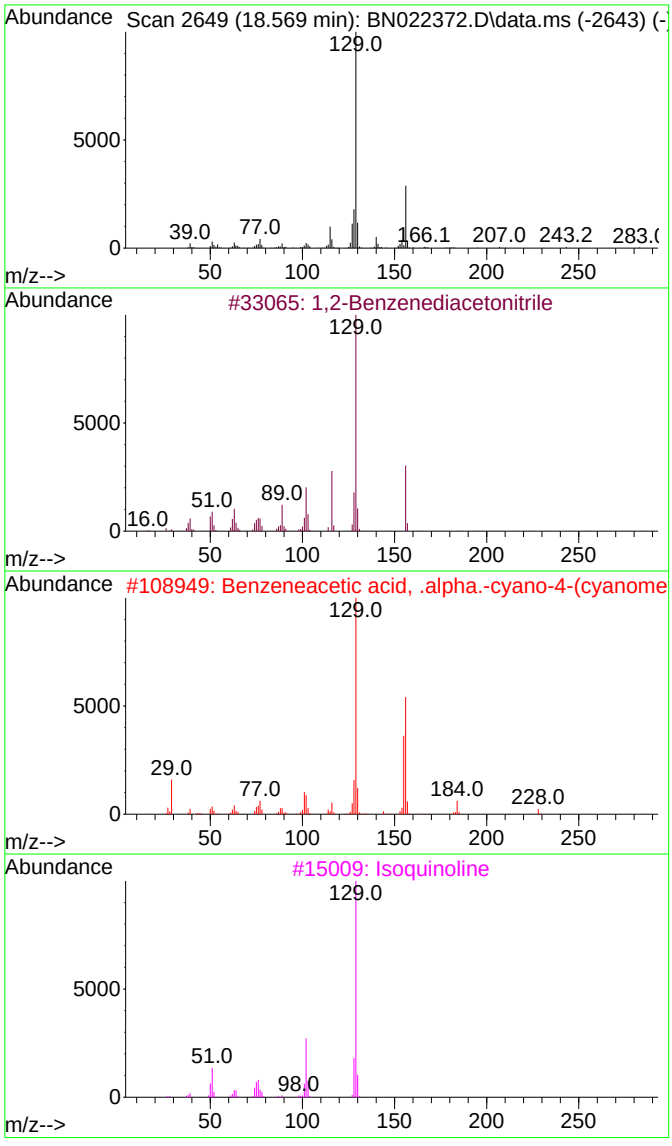
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

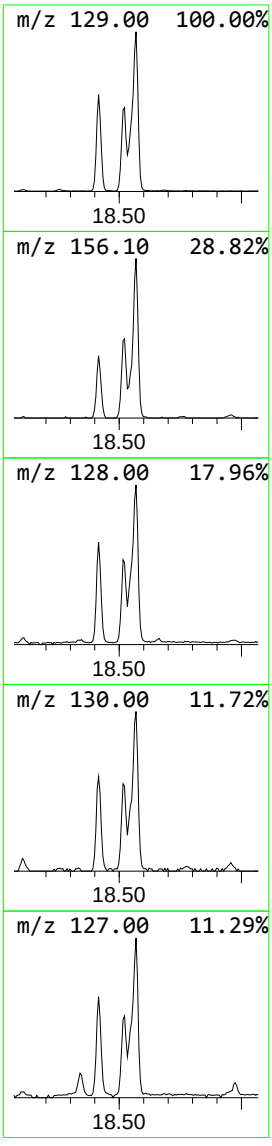
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1,2-Benzenediacetonitrile Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.569	5.13 ng/ul	450979	Phenanthrene-d10	17.363

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	59
2	Benzeneacetic acid, .alpha.-cyan...	228	C13H12N2O2	134882-04-5	50
3	Isoquinoline	129	C9H7N	000119-65-3	49
4	1-Deuteronaphthalene	129	C10H7D	000875-62-7	47
5	Benzeneacetonitrile, .alpha.-met...	129	C9H7N	000495-10-3	47





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Sample : N5232-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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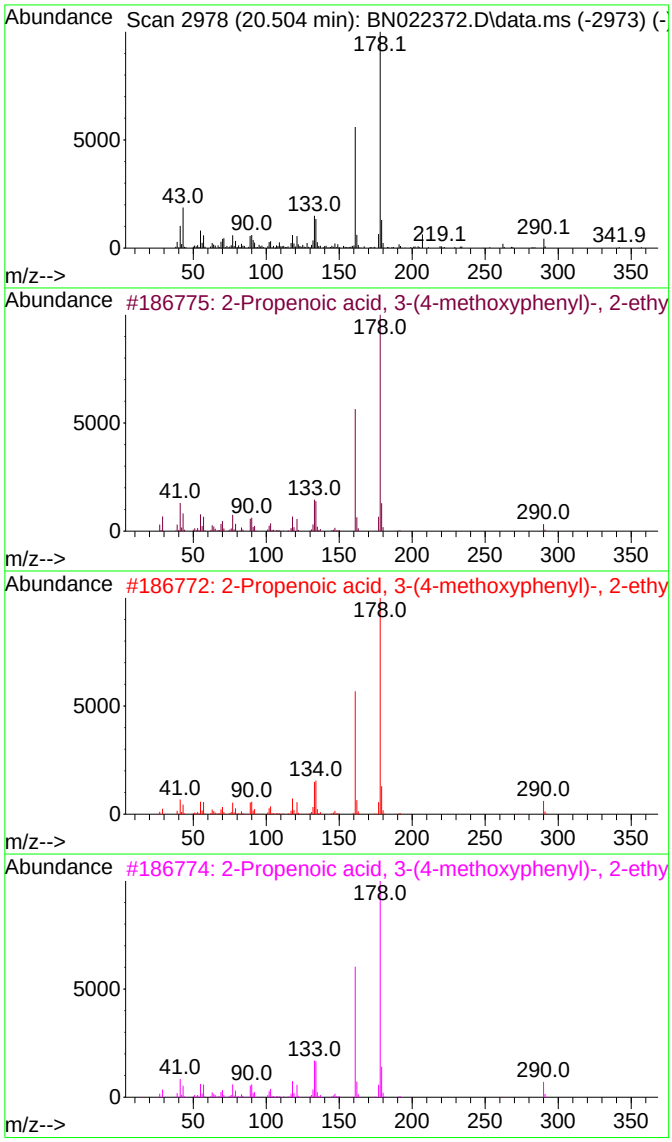
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

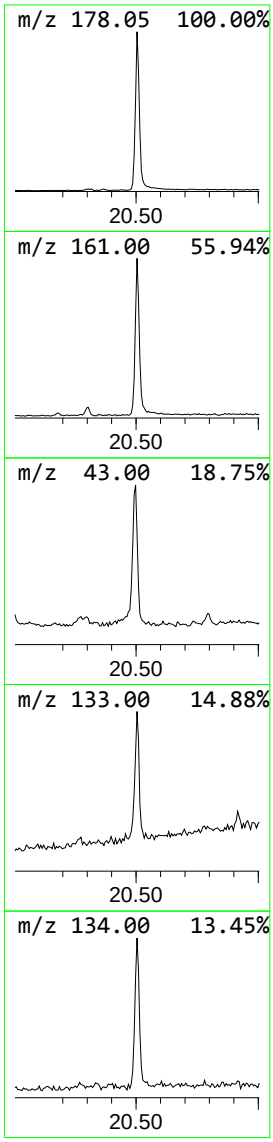
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2-Propenoic acid, 3-(4-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.504	2.83 ng/ul	250006	Chrysene-d12	21.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	99
2	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	99
3	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	97
4	2-Ethylhexyl trans-4-methoxycinn...	290	C18H26O3	083834-59-7	94
5	2-Propenoic acid, 3-(4-methoxyph...	290	C18H26O3	005466-77-3	91





Tentatively Identified Compound (LSC) summary

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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	3.034	5.0	ng/ul	282349	1	7.934	1121480	20.0
Ethanol, 1-(2-b...	10.534	6.9	ng/ul	481455	2	10.763	1392410	20.0
Benzenamine, 2,...	10.851	3.0	ng/ul	210734	2	10.763	1392410	20.0
Phenol, p-tert-...	12.110	6.7	ng/ul	462771	2	10.763	1392410	20.0
4-Hydroxy-N,N-d...	18.039	3.2	ng/ul	277608	4	17.363	1757270	20.0
Diphenyl sulfone	18.339	2.1	ng/ul	181042	4	17.363	1757270	20.0
Quinoline	18.416	2.7	ng/ul	237476	4	17.363	1757270	20.0
Quinolinium, 1-...	18.516	2.2	ng/ul	194167	4	17.363	1757270	20.0
1,2-Benzenediac...	18.569	5.1	ng/ul	450979	4	17.363	1757270	20.0
2-Propenoic aci...	20.504	2.8	ng/ul	250006	5	21.563	1763790	20.0