LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

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-BP093022.M
Title : SVOA CALIBRATION

Signal : TIC: BP011992.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	3	8	14	rVB	4853934	5828361	22.16%	0.890%
2	4.046	174	180	186	rBV	1713428	2294529	8.72%	0.350%
3	5.369	399	405	422	rBV	3998188	6518532	24.78%	0.995%
4	5.505	422	428	445	rVV	1306083	3068967	11.67%	0.469%
5	5.875	481	491	509	rVB	1752618	2899905	11.03%	0.443%
6	7.516	765	770	777	rVB	1446380	2291287	8.71%	0.350%
7	7.593	777	783	796	rVB	2421519	3994051	15.19%	0.610%
8	7.987	841	850	858	rVB	1236086	2140348	8.14%	0.327%
9	8.252	888	895	902	rVV	13175702	24184292	91.95%	3.692%
10	8.322	902	907	913	rVV	3423101	5476254	20.82%	0.836%
11	8.416	916	923	930	rVB	7757974	12909001	49.08%	1.971%
12	8.781	980	985	994	rVV	1224145	2976054	11.31%	0.454%
13	8.869	994	1000	1009	rVV	2493718	4382869	16.66%	0.669%
14	9.040	1023	1029	1044	rVB3	1116805	2867170	10.90%	0.438%
15	9.234	1055	1062	1067	rVV	1469726	2323538	8.83%	0.355%
16	9.316	1067	1076	1080	rVV	10912055	21033908	79.97%	3.211%
17	9.363	1080	1084	1090	rVV	2897887	5391598	20.50%	0.823%
18	9.863	1162	1169	1172	rVV	5371799	10279372	39.08%	1.569%
19	9.899	1172	1175	1190	rVB2	7919766	14337012	54.51%	2.189%
20	10.081	1199	1206	1215	rBV4	1353243	4072530	15.48%	0.622%
21	10.181	1215	1223	1233	rVV	10069954	19654141	74.72%	3.001%
22	10.328	1239	1248	1254	rVV2	3322794	7307020	27.78%	1.116%
23	10.510	1274	1279	1283	rVV	1578001	2556687	9.72%	0.390%
24	10.563	1283	1288	1293	rVV	4115314	7200176	27.37%	1.099%
25	10.740	1312	1318	1326	rVV	9305885	17778397	67.59%	2.714%
26	10.846	1326	1336	1347	rVV2	11677329	26302029	100.00%	4.016%
27	11.046	1363	1370	1376	rBV	1289222	2648771	10.07%	0.404%
28	11.140	1382	1386	1395	rVV	1431356	2936958	11.17%	0.448%
29	11.228	1395	1401	1405	rVV	2072324	3624926	13.78%	0.553%
30	11.281	1405	1410	1430	rVB	4387659	11089347	42.16%	1.693%
31	11.534	1442	1453	1463	rVV	6385312	13255468	50.40%	2.024%
32	11.722	1477	1485	1489	rBV	7977973	13652396	51.91%	2.084%
33	11.775	1489	1494	1498	rVV	5476709	9330901	35.48%	1.425%
34	11.816	1498	1501	1508	rVV2	1949242	3322777	12.63%	0.507%
35	12.093	1543	1548	1555	rVB3	1314076	2417156	9.19%	0.369%
36	12.198	1561	1566	1570	rBV2	1443318	2607221	9.91%	0.398%

LSC Area Percent Report

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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

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-BP093022.M
Title : SVOA CALIBRATION

37	12.393	1594	1599	1604	rBV	2310676	3933398	14.95%	0.601%
38	12.451	1604	1609	1618	rVB2	4274418	7310077	27.79%	1.116%
39	12.569	1618	1629	1635	rBV	11206304	20499942	77.94%	3.130%
40	12.675	1643	1647	1651	rVB	2115217	2842848	10.81%	0.434%
41	12.722	1651	1655	1662	rBV3	1306639	2902662	11.04%	0.443%
42	12.945	1687	1693	1698	rBV3	1519067	2619517	9.96%	0.400%
43	13.040	1705	1709	1718	rVB4	1045512	2200043	8.36%	0.336%
44	13.322	1752	1757	1759	rBV	1464427	2562163	9.74%	0.391%
45	13.351	1759	1762	1772	rVB2	2822802	5886930	22.38%	0.899%
46	13.663	1808	1815	1820	rBV3	2655946	5603654	21.31%	0.856%
47	13.845	1840	1846	1852	rVV3	3236778	6925175	26.33%	1.057%
48	13.940	1858	1862	1866	rVV	1616691	2431455	9.24%	0.371%
49	14.016	1870	1875	1880	rVV	2140208	3032600	11.53%	0.463%
50	14.122	1889	1893	1899	rVV3	1126742	2394028	9.10%	0.366%
51	14.216	1904	1909	1911	rVV	2407369	3807800	14.48%	0.581%
52	14.245	1911	1914	1921	rVB	3467669	5443411	20.70%	0.831%
53	14.522	1956	1961	1967	rVB	1432349	2229352	8.48%	0.340%
54	14.587	1967	1972	1979	rVB	8188978	13065252	49.67%	1.995%
55	14.657	1979	1984	1988	rBV	1481021	2267276	8.62%	0.346%
56	14.710	1988	1993	2003	rVB5	1387893	3281490	12.48%	0.501%
57	14.810	2004	2010	2017	rBV3	2122929	4230090	16.08%	0.646%
58	14.922	2025	2029	2036	rVB2	4980518	7932541	30.16%	1.211%
59	15.122	2045	2063	2071	rVV	5484447	19770436	75.17%	3.018%
60	15.457	2112	2120	2125	rVV	4819814	10602305	40.31%	1.619%
61	15.516	2125	2130	2135	rVV	2945804	5172061	19.66%	0.790%
62	15.575	2135	2140	2147	rVV	3342567	5861666	22.29%	0.895%
63	15.639	2147	2151	2157	rVV6	1258995	2500189	9.51%	0.382%
64	15.716	2157	2164	2171	rVV	4842075	11189016	42.54%	1.708%
65	15.810	2171	2180	2186	rVV3	3108119	9227219	35.08%	1.409%
66	15.934	2194	2201	2203	rVV	3362967	5860065	22.28%	0.895%
67	15.963	2203	2206	2210	rVV	2993460	5092852	19.36%	0.778%
68	16.004	2210	2213	2220	rVV3	2511939	4356242	16.56%	0.665%
69	16.075	2220	2225	2234	rVB4	934685	2060185	7.83%	0.315%
70	16.304	2261	2264	2270	rVB2	5730454	10370992	39.43%	1.583%
71	16.492	2288	2296	2299	rBV3	2550500	5324749	20.24%	0.813%
72	16.551	2299	2306	2308	rBV2	3084979	5526867	21.01%	0.844%
73	16.928	2354	2370	2375	rVB2	3024086	10551055	40.11%	1.611%
74	17.063	2380	2393	2402	rBV3	3540140	10173629	38.68%	1.553%
75	17.275	2415	2429	2434	rVV3	7421148	25296079	96.18%	3.862%
76	17.328	2434	2438	2444	rVB	6877024	9703394	36.89%	1.481%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

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-BP093022.M
Title : SVOA CALIBRATION

77	17.381	2444	2447	2451	rVB2	2281464	2607156	9.91%	0.398%
78	17.480	2457	2464	2470	rBV4	2790076	7692360	29.25%	1.174%
79	17.581	2477	2481	2486	rVB	1272837	2090463	7.95%	0.319%
80	17.663	2486	2495	2497	rBV2	1500507	3474898	13.21%	0.531%
81	17.698	2497	2501	2505	rBV	1147014	2203298	8.38%	0.336%
82	17.792	2512	2517	2523	rVB4	1762918	3364664	12.79%	0.514%
83	17.857	2523	2528	2536	rBV2	2178276	4030060	15.32%	0.615%
84	17.998	2547	2552	2556	rBV2	1128383	2043859	7.77%	0.312%
85	18.122	2568	2573	2580	rBV3	2821118	4740214	18.02%	0.724%
86	18.192	2580	2585	2592	rBV6	1814691	4655189	17.70%	0.711%
87	18.339	2605	2610	2620	rVB3	2173839	4219269	16.04%	0.644%
88	18.428	2620	2625	2628	rBV	1455363	2564778	9.75%	0.392%
89	19.051	2725	2731	2737	rBV3	1157248	2336772	8.88%	0.357%
90	19.292	2767	2772	2778	rVB	5461862	7187139	27.33%	1.097%
91	19.398	2784	2790	2793	rBV3	1198156	2076290	7.89%	0.317%
92	19.616	2822	2827	2830	rBV2	5109995	8072702	30.69%	1.232%
93	19.645	2830	2832	2840	rVV2	4502904	6724153	25.57%	1.027%
94	20.098	2902	2909	2912	rBV	1958323	3541955	13.47%	0.541%
95	21.363	3118	3124	3128	rBV2	3336931	6717543	25.54%	1.026%
96	21.404	3128	3131	3142	rVB	1800653	2630553	10.00%	0.402%
97	23.051	3406	3411	3417	rBV	1461803	3542121	13.47%	0.541%
98	23.639	3505	3511	3516	rVV2	2909108	5607537	21.32%	0.856%
99	23.686	3516	3519	3528	rVB	1382066	2412494	9.17%	0.368%
100	23.798	3532	3538	3543	rBV2	1722161	3459137	13.15%	0.528%

Sum of corrected areas: 654989258

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
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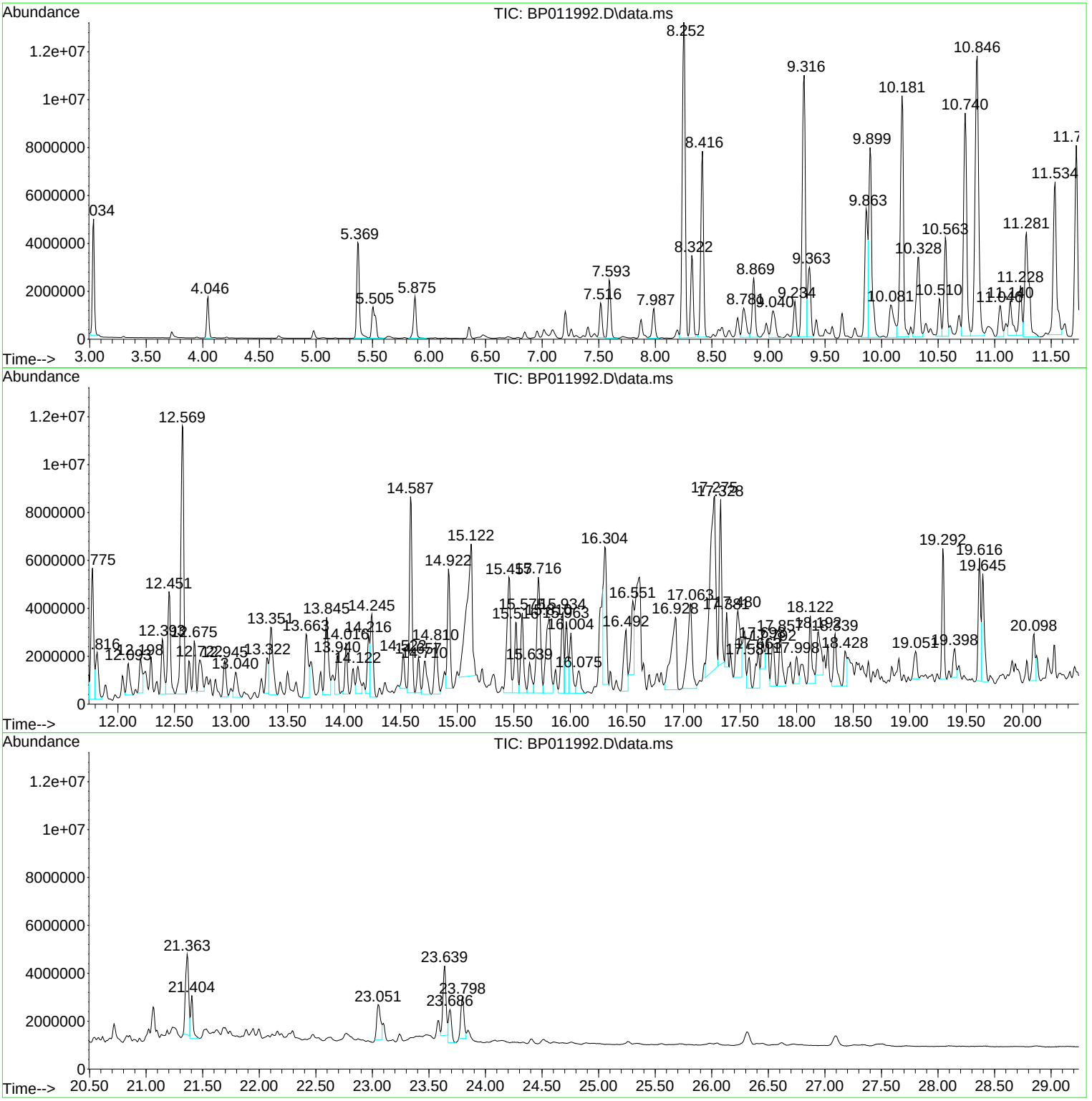
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
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Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

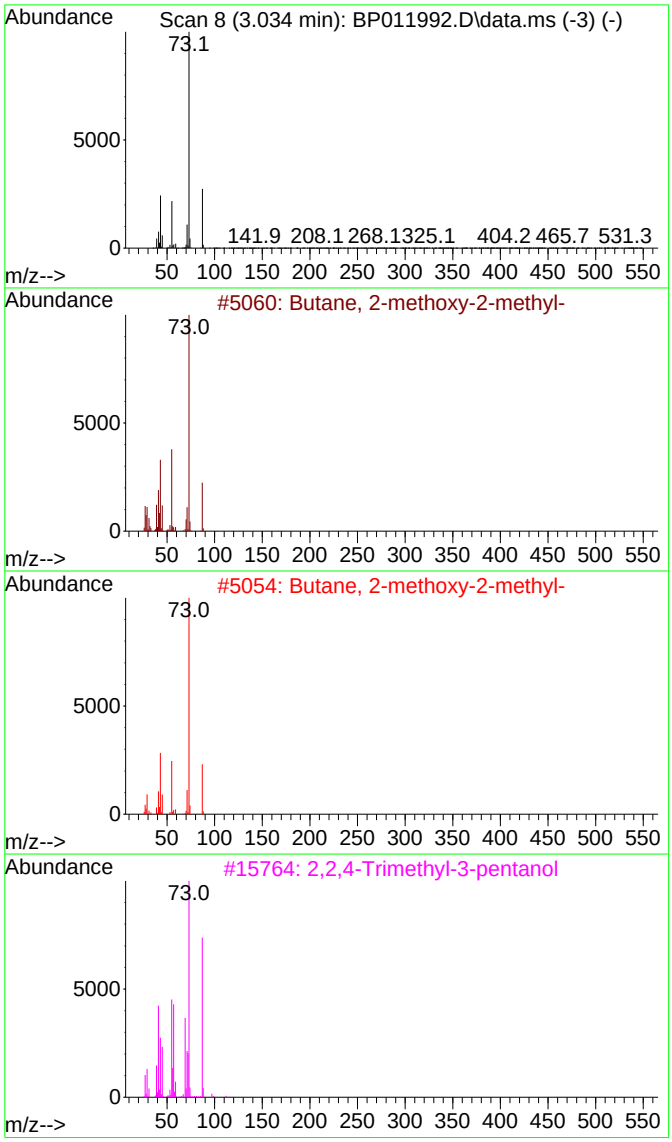
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

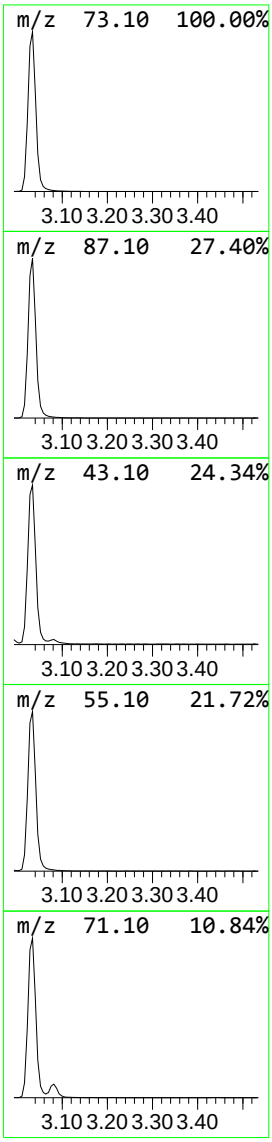
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	54.46 ng/ul	5828360	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

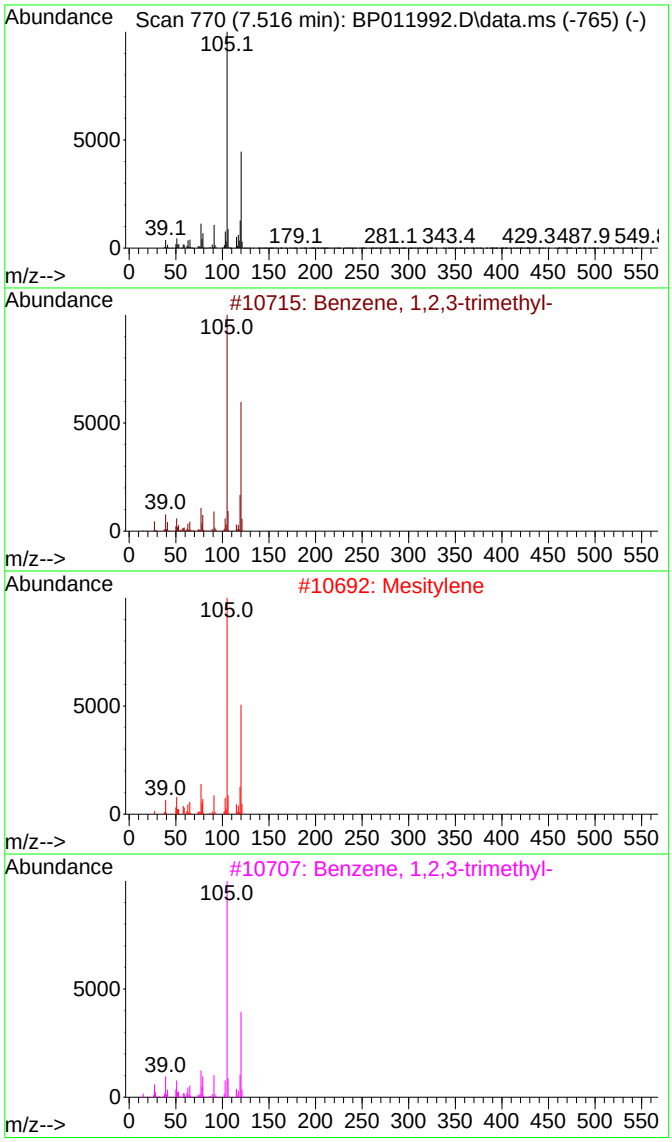
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

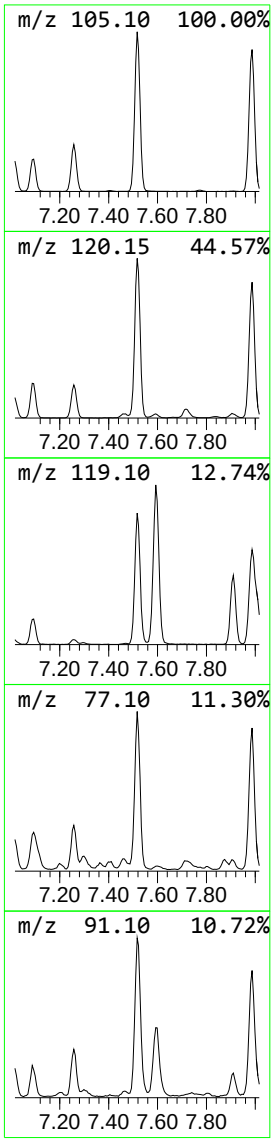
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.516	21.41 ng/ul	2291290	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
2	Mesitylene	120	C9H12	000108-67-8	93
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

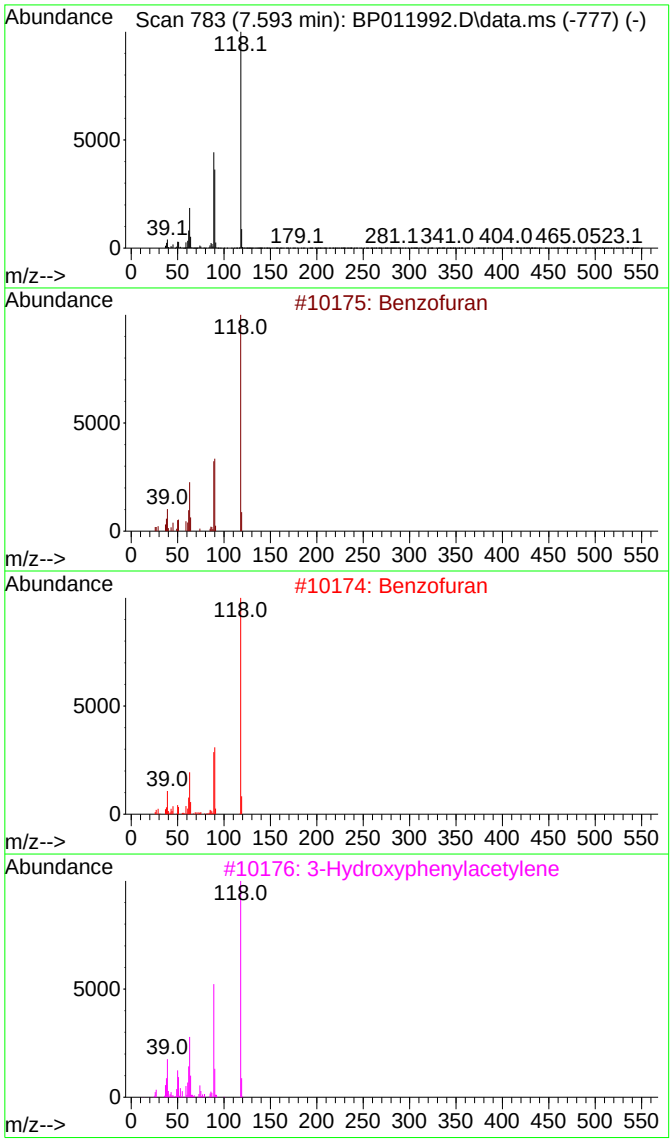
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

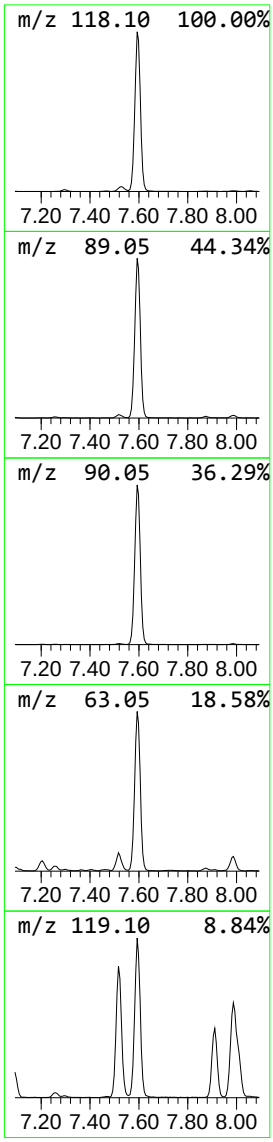
****

Peak Number 7 Benzofuran Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.593	37.32 ng/ul	3994050	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	95
2	Benzofuran	118	C8H6O	000271-89-6	91
3	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	90
4	Benzofuran	118	C8H6O	000271-89-6	87
5	Benzofuran	118	C8H6O	000271-89-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

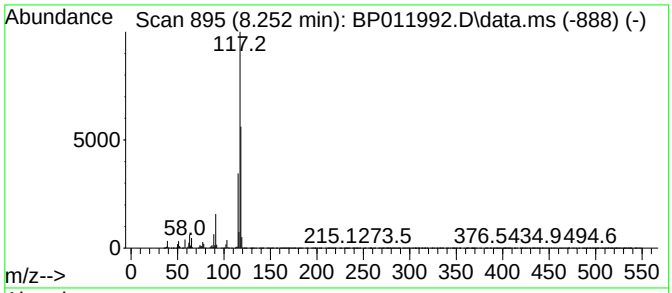
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

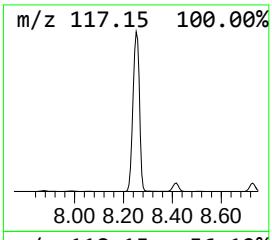
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

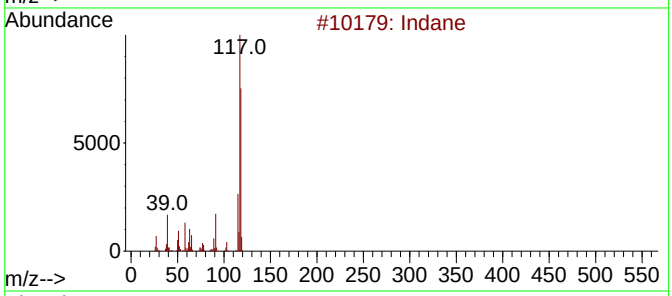
****
Peak Number 9 Indane Concentration Rank 1

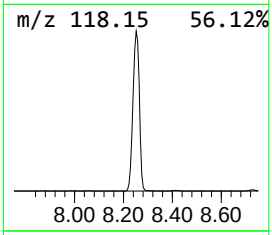
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.252	225.99 ng/ul	24184300	1,4-Dichlorobenzene-d4	7.875

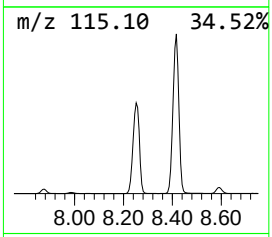
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	Indane	118	C9H10	000496-11-7	87
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	83
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
5	Deltacyclene	118	C9H10	007785-10-6	72

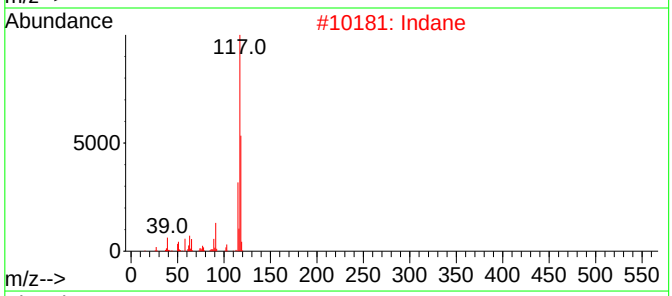


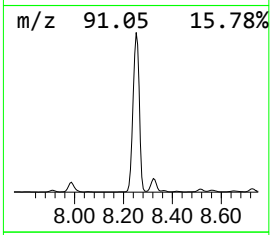


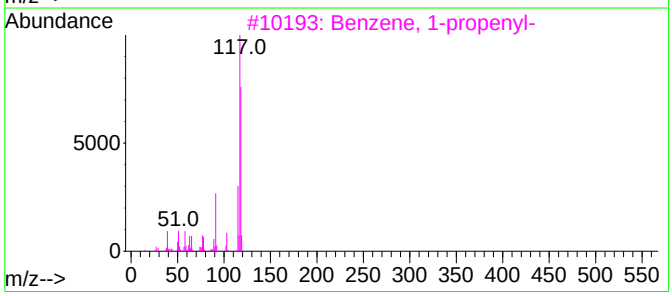


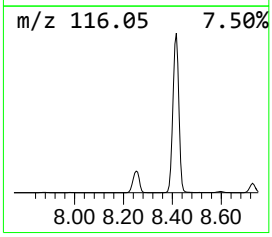












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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

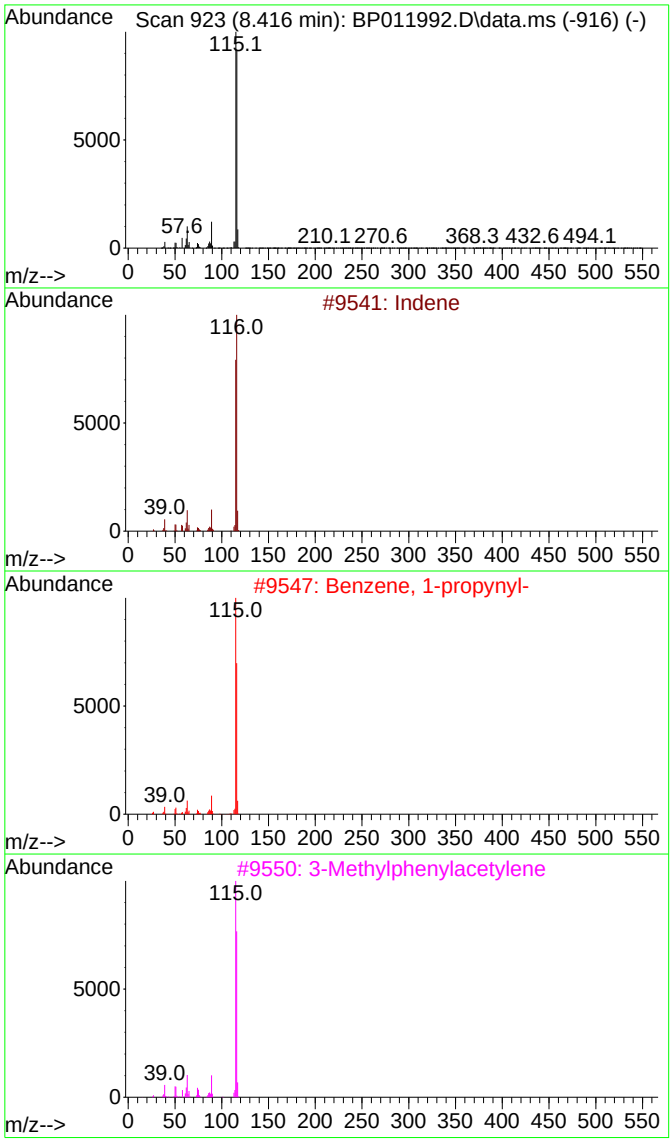
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

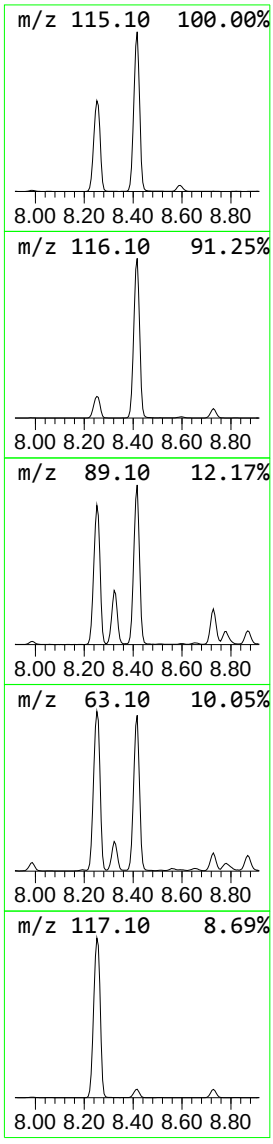
****

Peak Number 10 Indene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.416	120.63 ng/ul	12909000	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

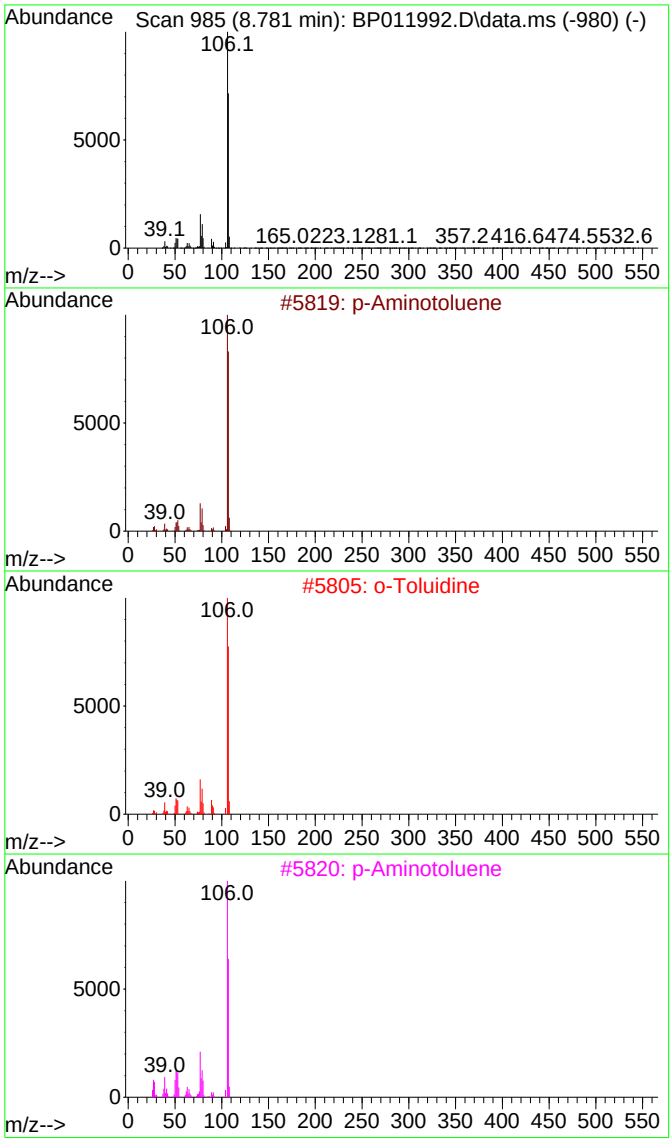
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

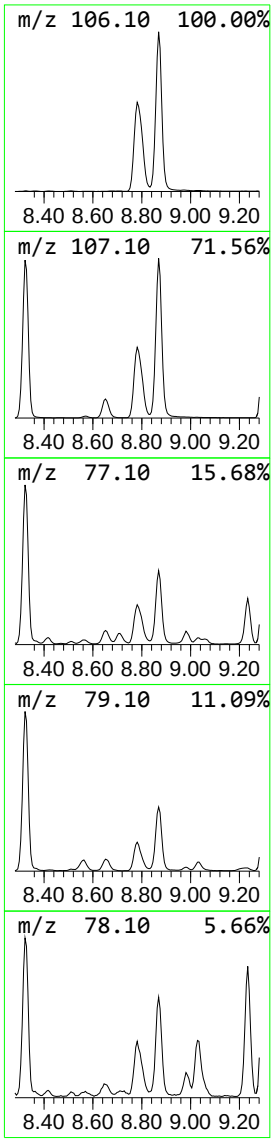
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 p-Aminotoluene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.781	27.81 ng/ul	2976050	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Aminotoluene	107	C7H9N	000106-49-0	94
2	o-Toluidine	107	C7H9N	000095-53-4	94
3	p-Aminotoluene	107	C7H9N	000106-49-0	91
4	o-Toluidine	107	C7H9N	000095-53-4	90
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	90





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

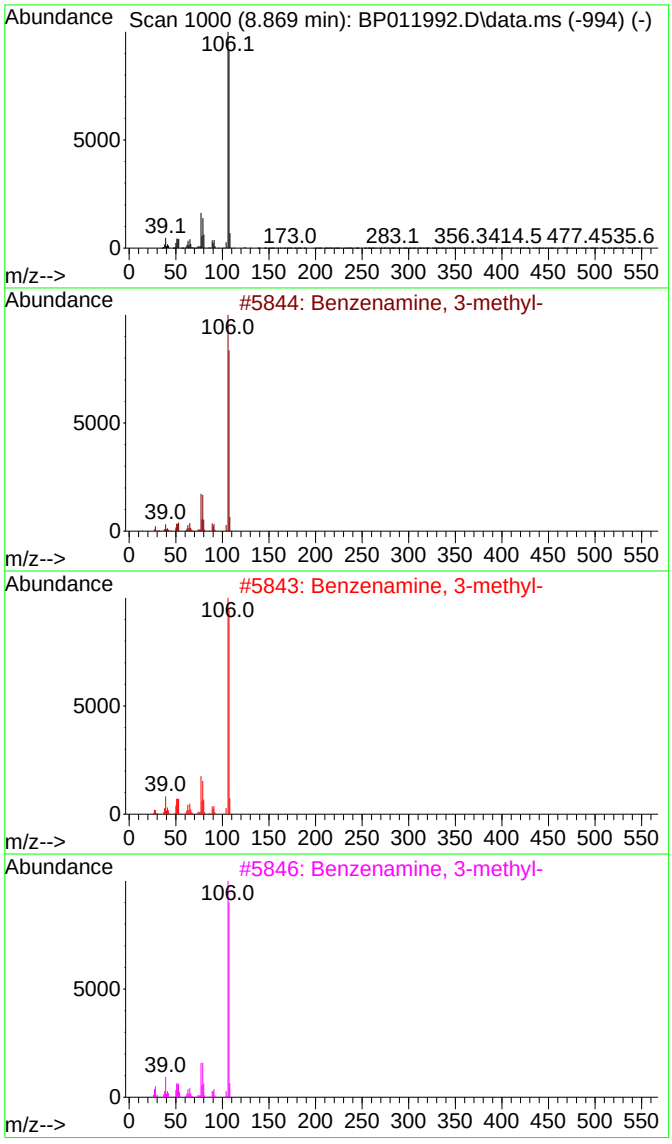
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

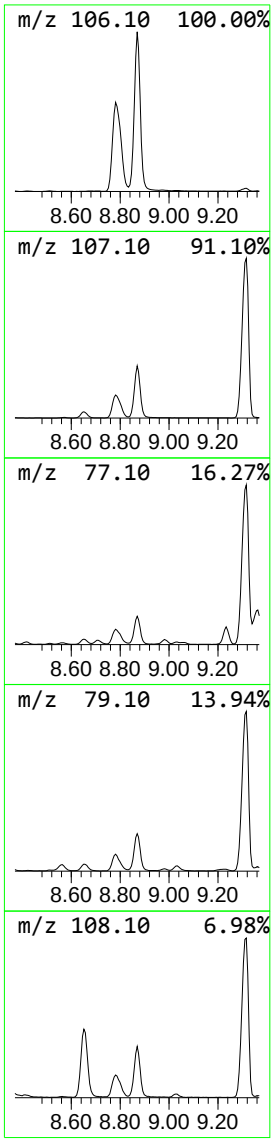
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzenamine, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.869	40.95 ng/ul	4382870	1,4-Dichlorobenzene-d4	7.875

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	95
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

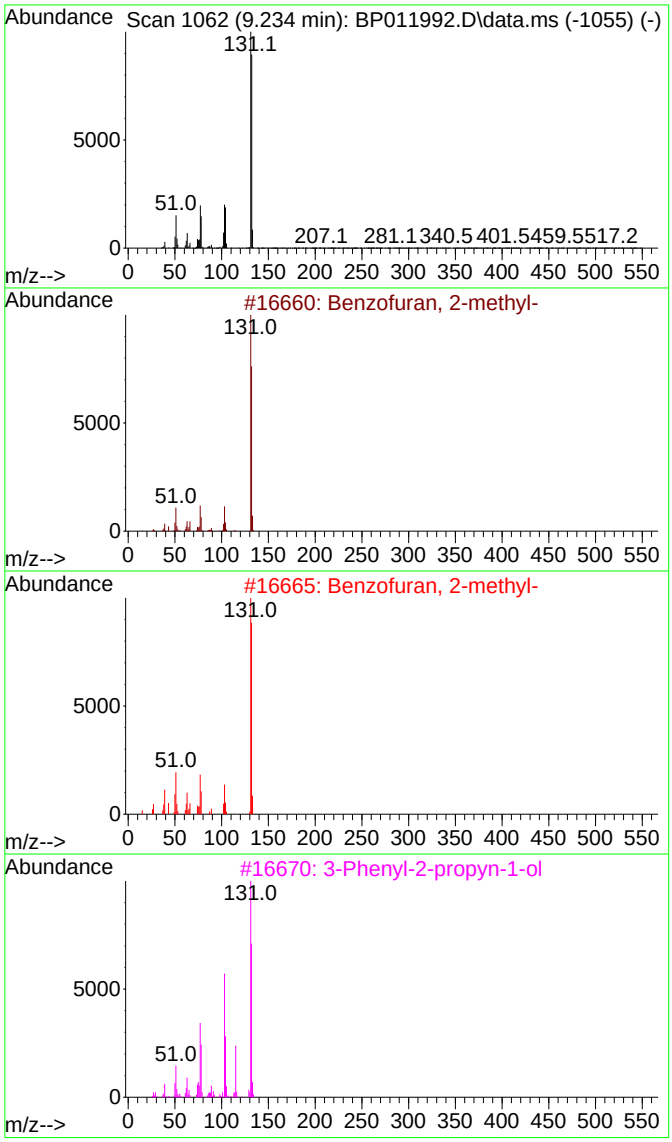
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

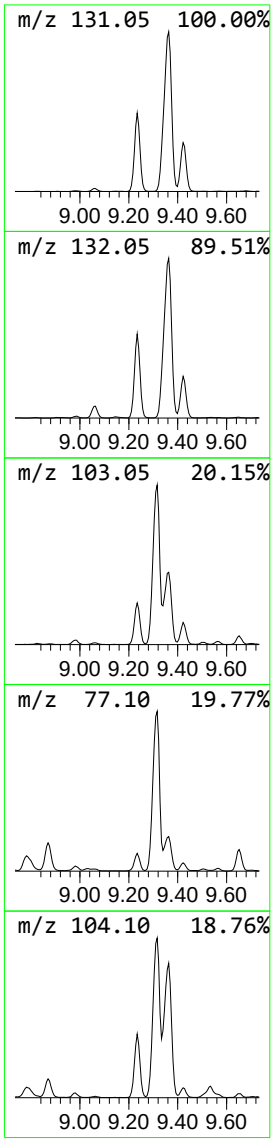
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzofuran, 2-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.234	21.71 ng/ul	2323540	1,4-Dichlorobenzene-d4	7.875

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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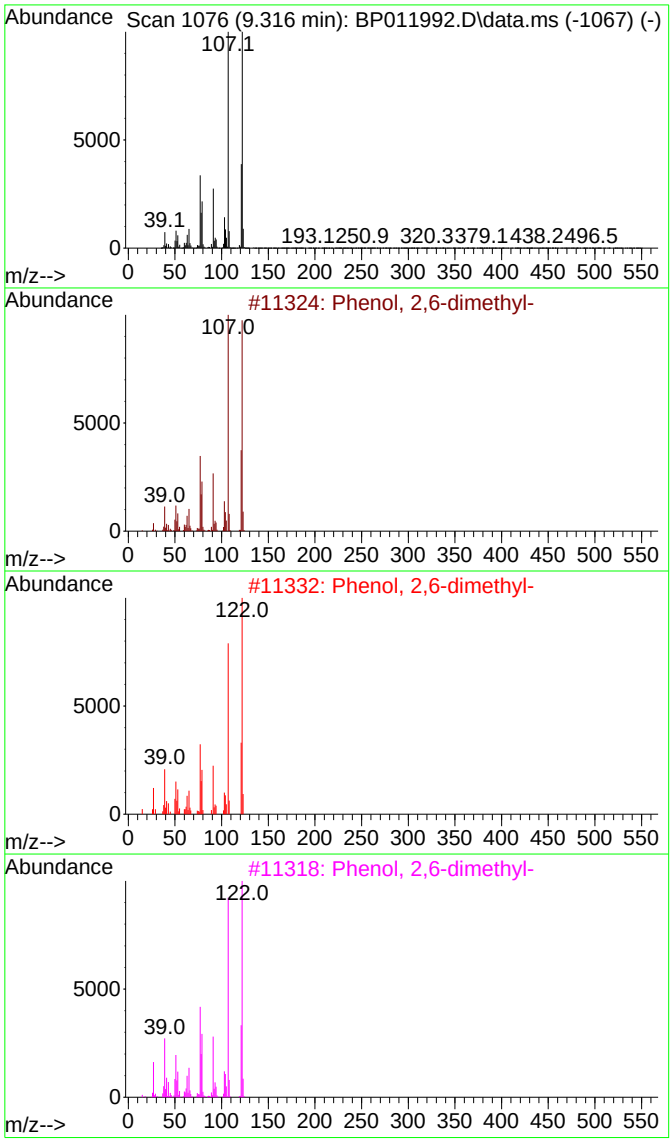
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

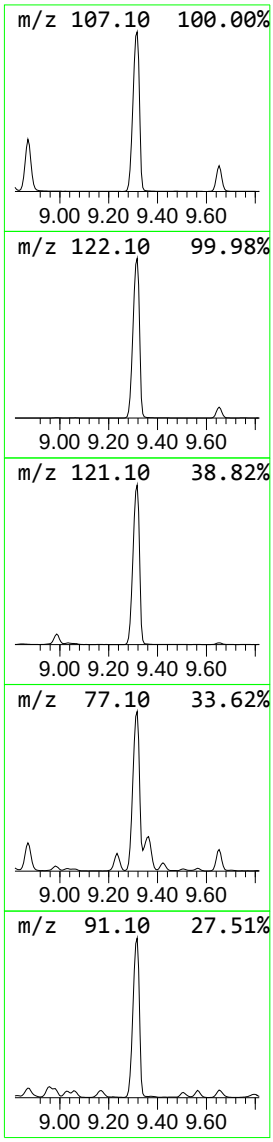
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Phenol, 2,6-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.316	23.66 ng/ul	21033900	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	97
2	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	96
3	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	96
4	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	94
5	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	94





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

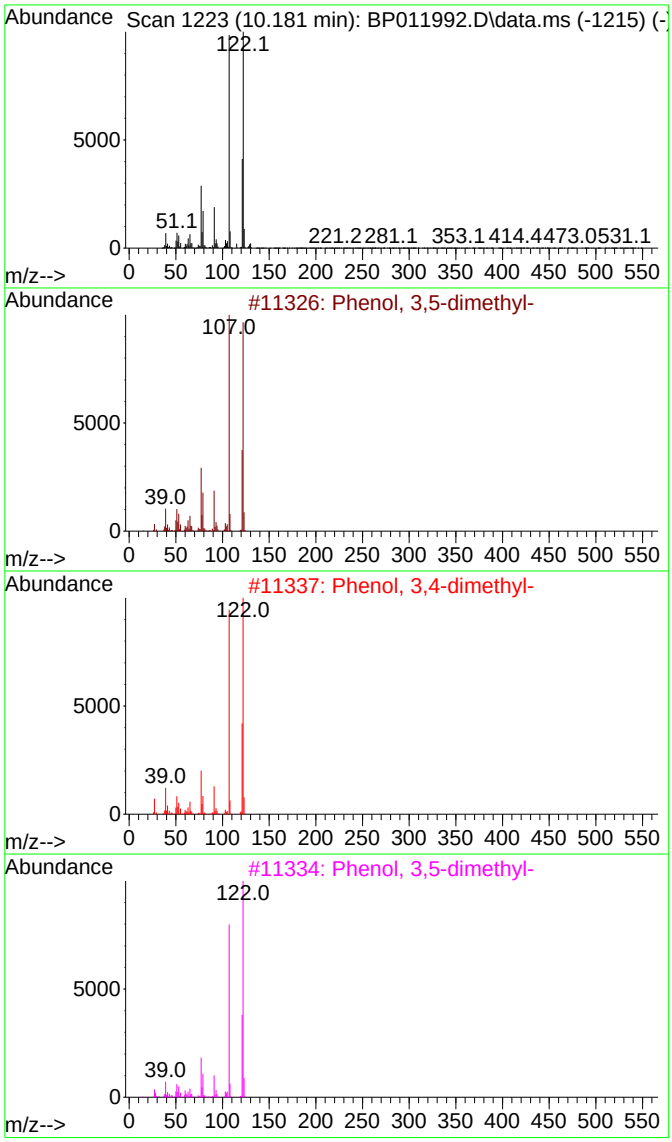
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

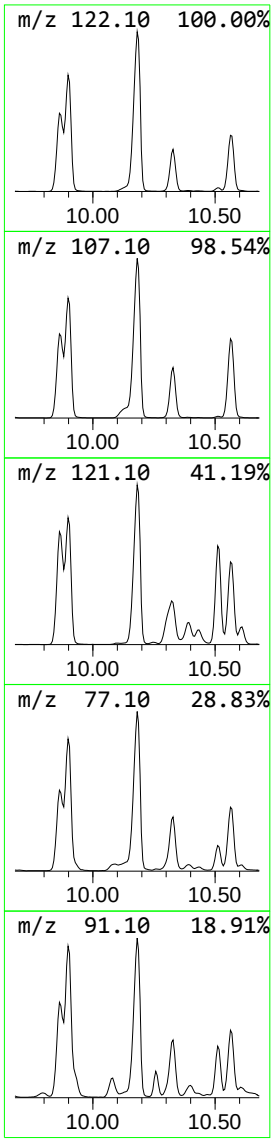
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Phenol, 3,5-dimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.181	22.11 ng/ul	19654100	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	97
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	97
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	96
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
5	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

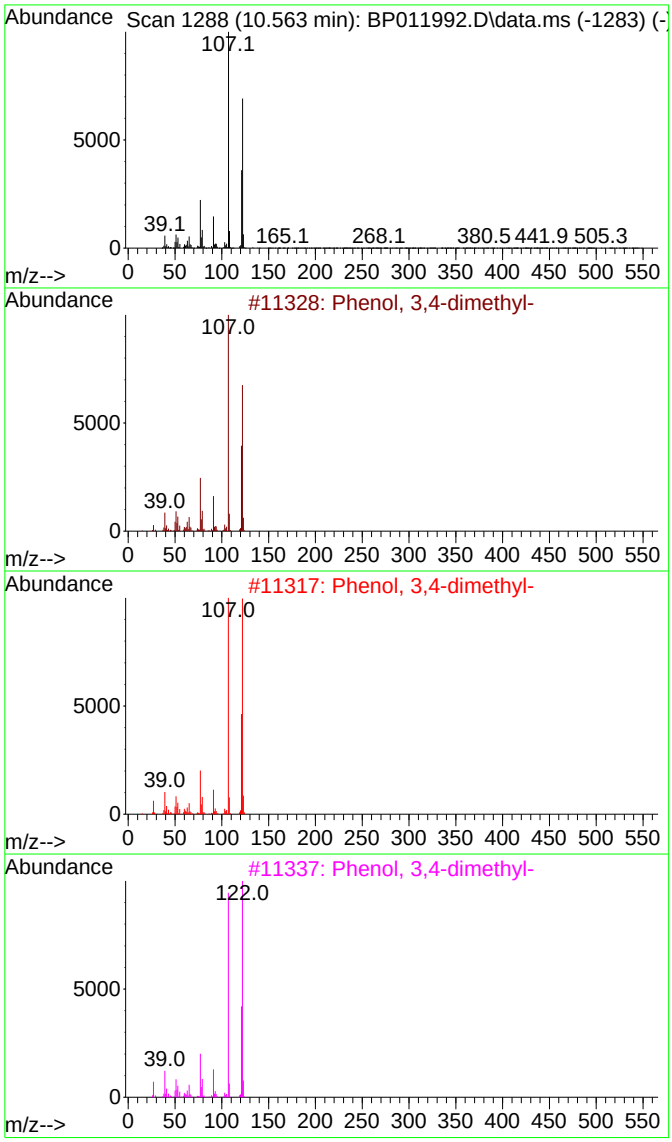
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

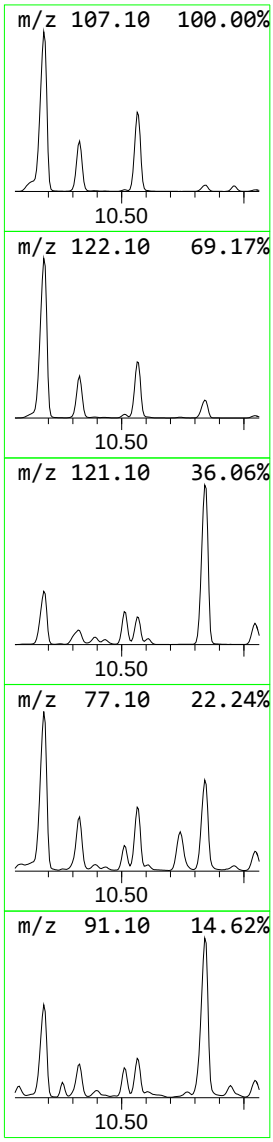
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Phenol, 3-ethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.563	8.10 ng/ul	7200180	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
5	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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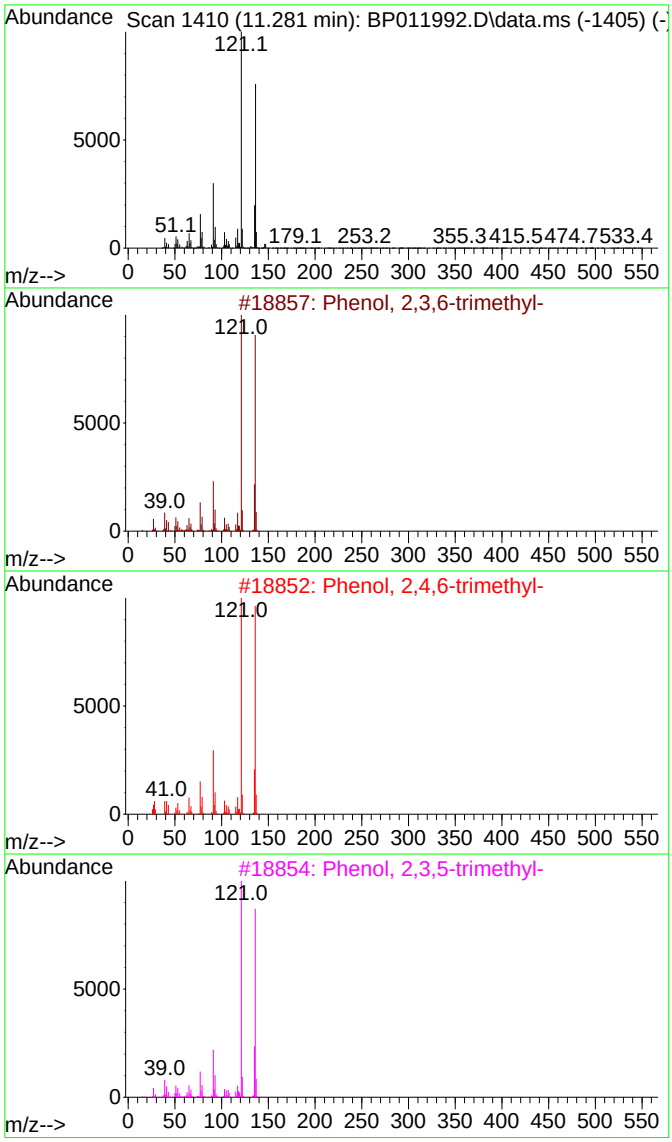
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

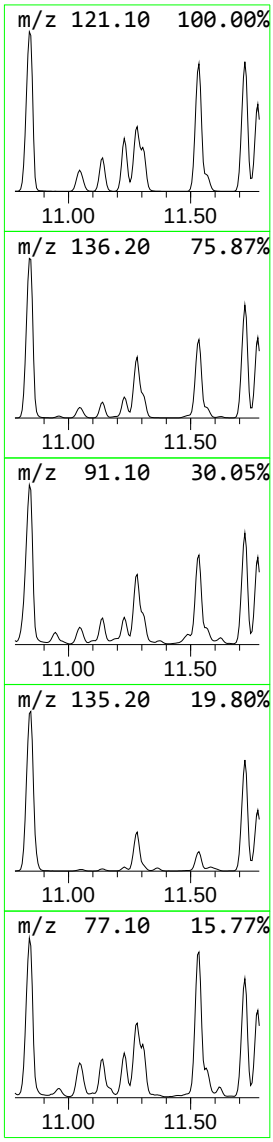
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Phenol, 2,3,6-trimethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.281	12.48 ng/ul	11089300	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	96
2	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	96
3	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	96
4	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95
5	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

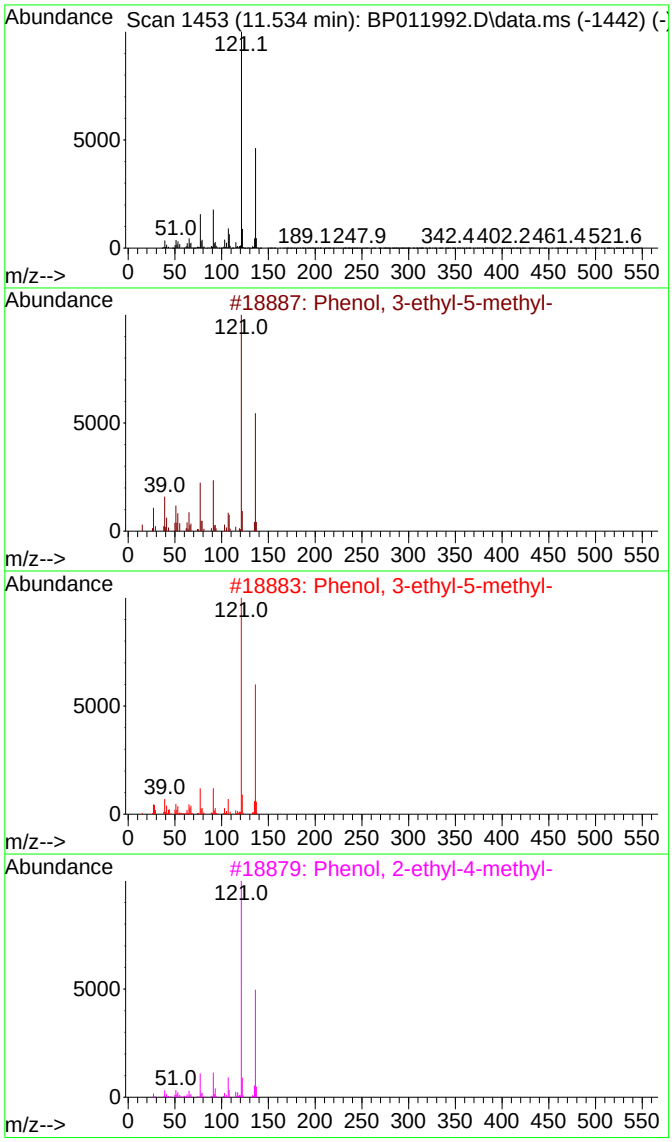
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

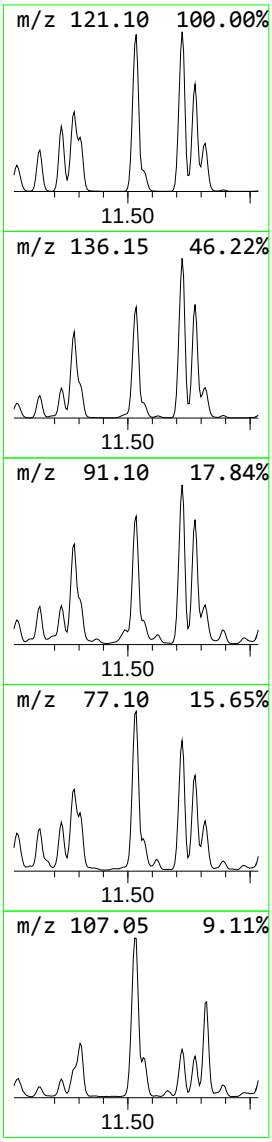
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Phenol, 3-ethyl-5-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.534	14.91 ng/ul	13255500	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	94
2	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	91
3	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	91
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	90
5	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	80





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Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

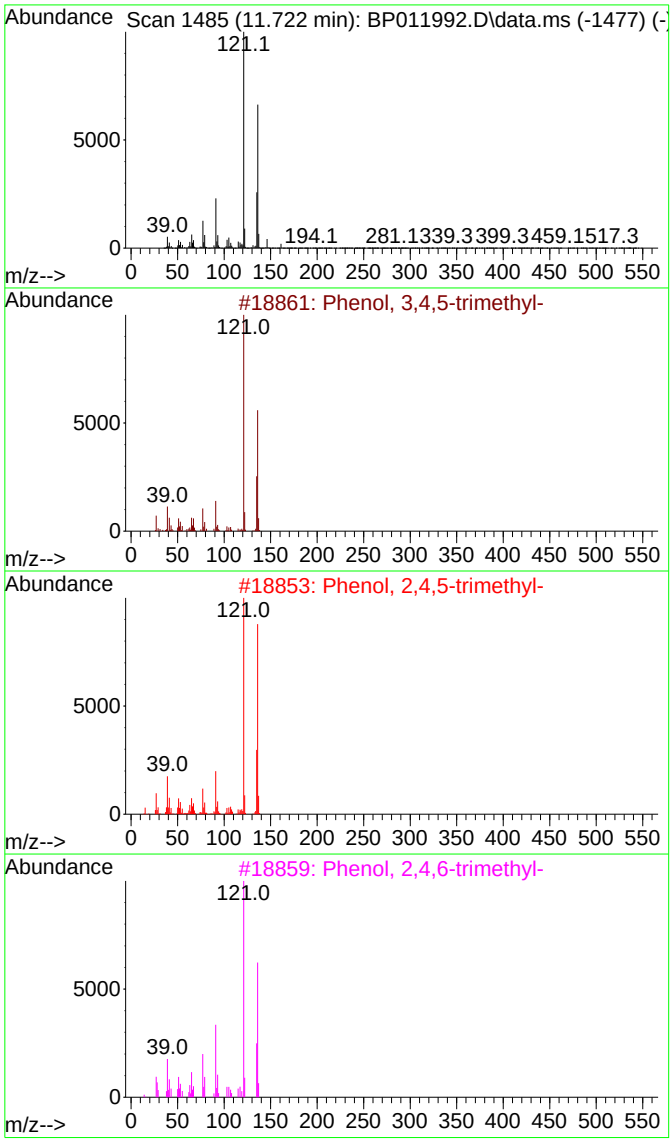
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

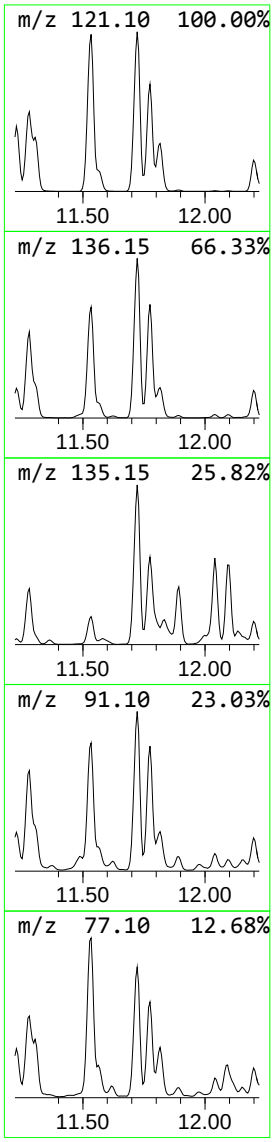
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Phenol, 3,4,5-trimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.722	15.36 ng/ul	13652400	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	94
2	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	94
3	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	91
4	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	91
5	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	91





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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

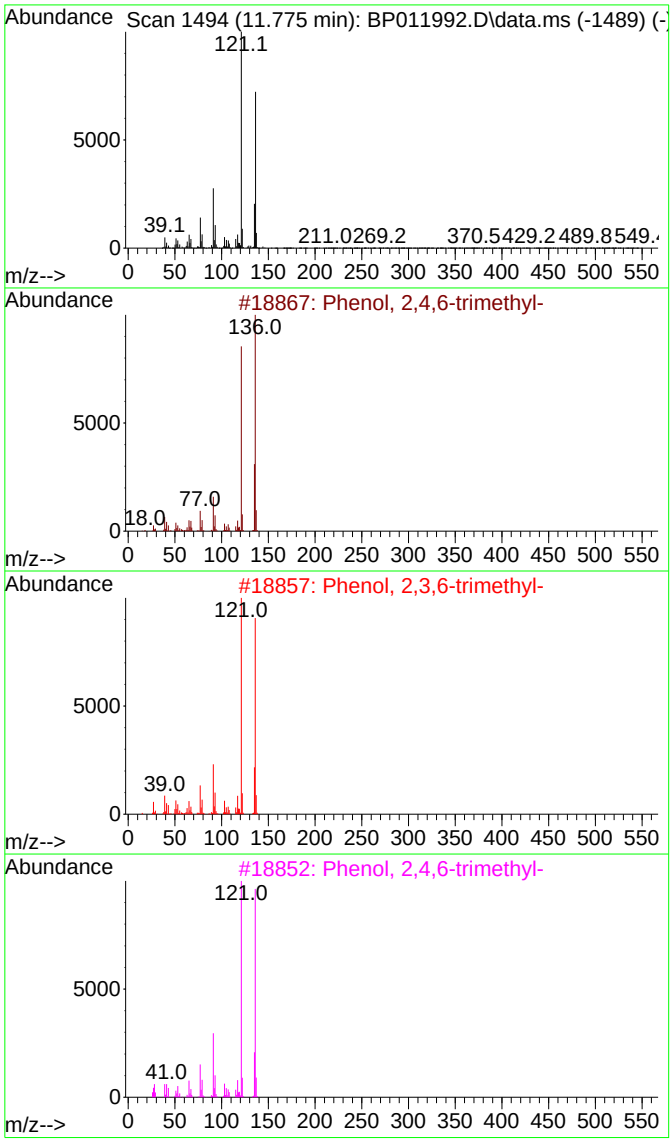
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

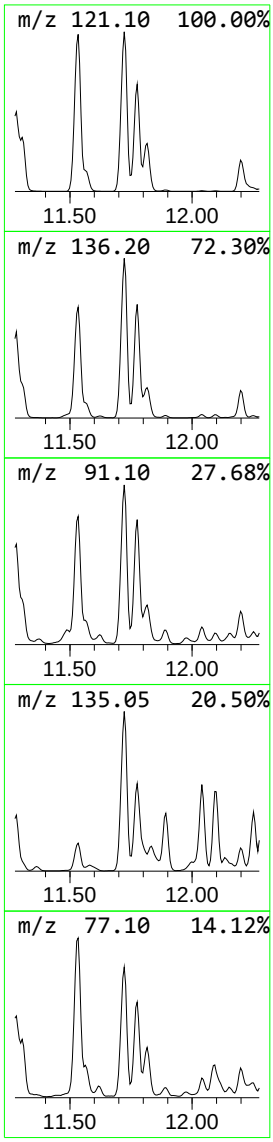
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Phenol, 2,4,6-trimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.775	10.50 ng/ul	9330900	Naphthalene-d8	10.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	97
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	96
3	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	96
4	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	96
5	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

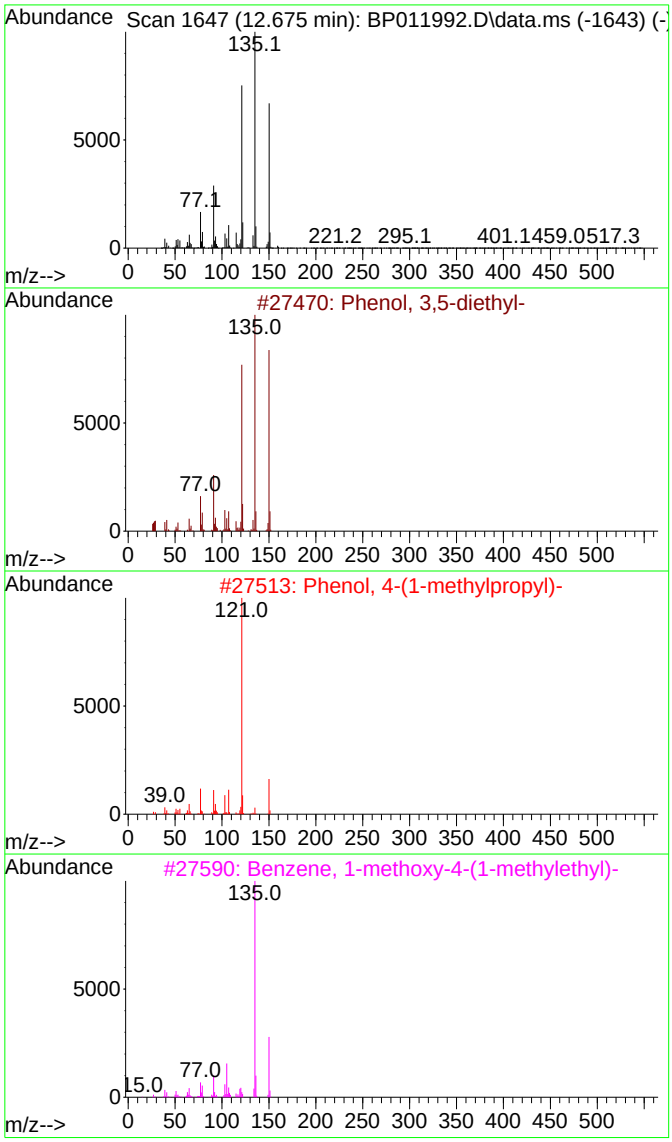
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

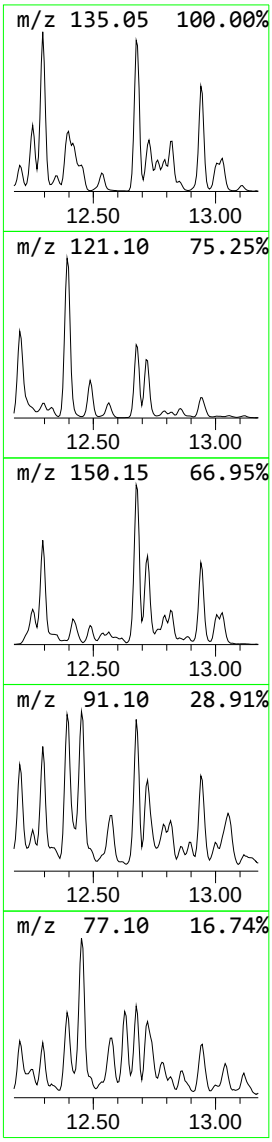
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 Phenol, 3,5-diethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.675	25.50 ng/ul	2842850	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-diethyl-	150	C10H14O	001197-34-8	96
2	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	53
3	Benzene, 1-methoxy-4-(1-methylet...	150	C10H14O	004132-48-3	53
4	2,3,5-Trimethylanizole	150	C10H14O	1000342-42-3	49
5	Phenol, 2,3,5,6-tetramethyl-	150	C10H14O	000527-35-5	49





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

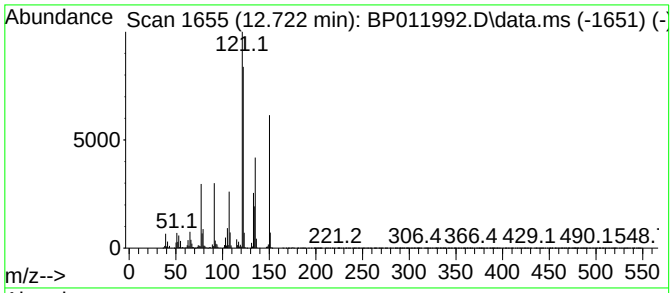
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

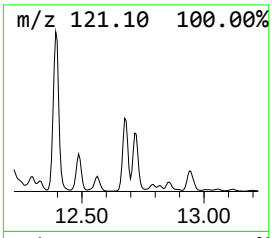
****

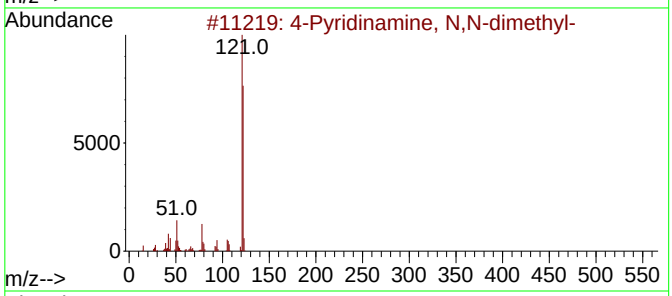
Peak Number 32 unknown-01 Concentration Rank 18

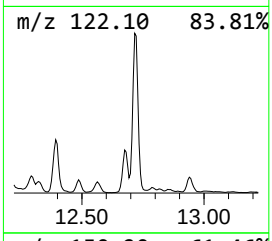
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.722	26.04 ng/ul	2902660	Acenaphthene-d10	14.522

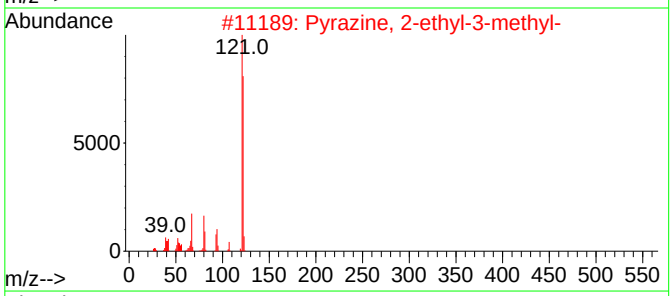
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Pyridinamine, N,N-dimethyl-	122	C7H10N2	001122-58-3	43
2	Pyrazine, 2-ethyl-3-methyl-	122	C7H10N2	015707-23-0	38
3	1(2H)-Naphthalenone, 3,4,5,6,7,8...	150	C10H14O	018631-96-4	38
4	2-sec-Butylphenol, O-isopropylox...	236	C14H20O3	1000487-30-6	35
5	1,3-Benzodioxole	122	C7H6O2	000274-09-9	35

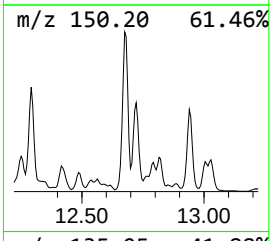


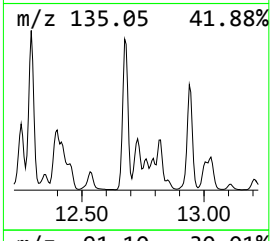


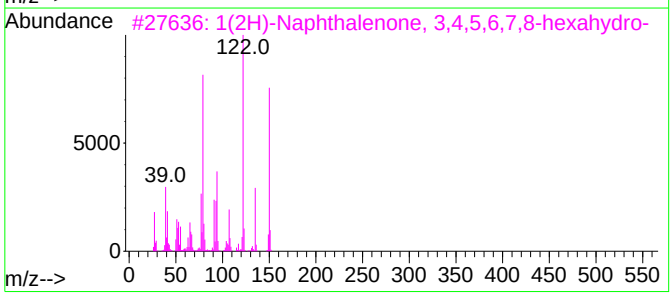


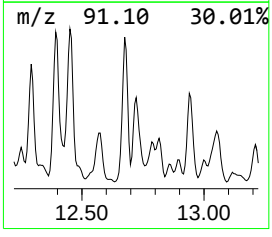












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

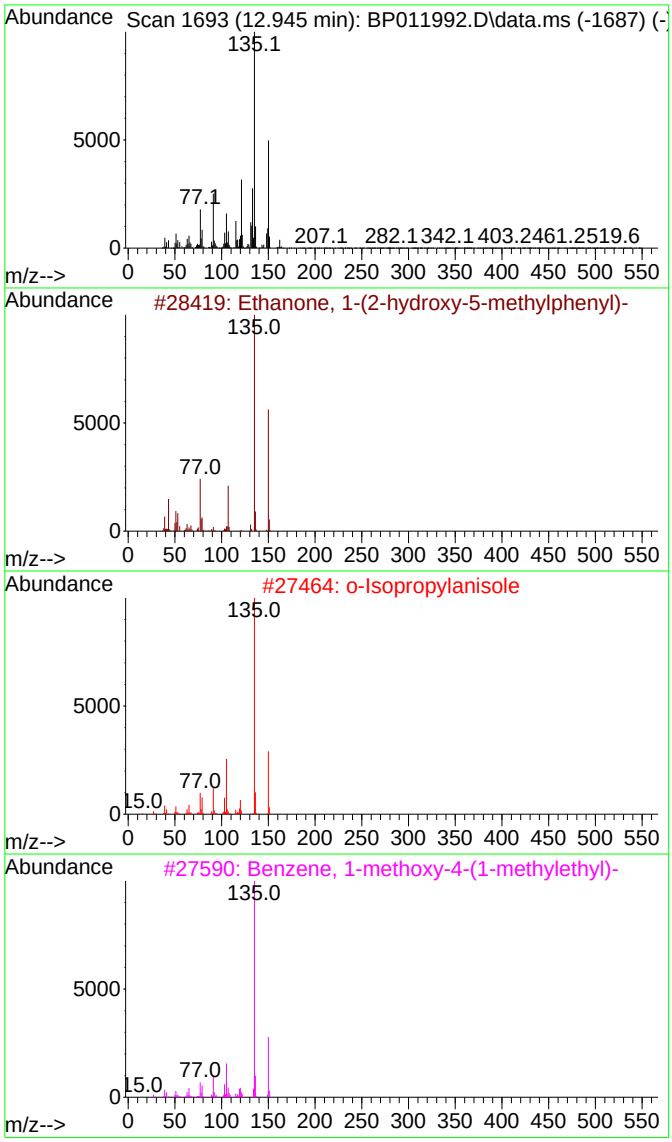
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

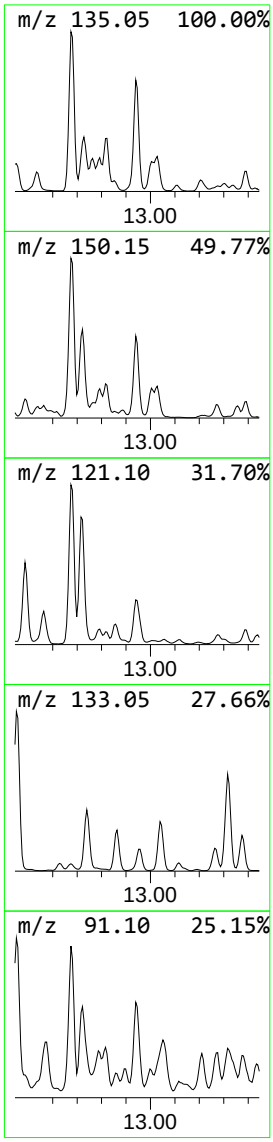
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 Ethanone, 1-(2-hydroxy-5-me... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.945	23.50 ng/ul	2619520	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(2-hydroxy-5-methylp...	150	C9H10O2	001450-72-2	70
2	o-Isopropylanisole	150	C10H14O	002944-47-0	70
3	Benzene, 1-methoxy-4-(1-methylet...	150	C10H14O	004132-48-3	70
4	3,6-Dimethylbenzene-1,2-diamine,...	150	C9H14N2	1000499-85-0	64
5	4-Hydroxy-2-methylacetophenone	150	C9H10O2	000875-59-2	62





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

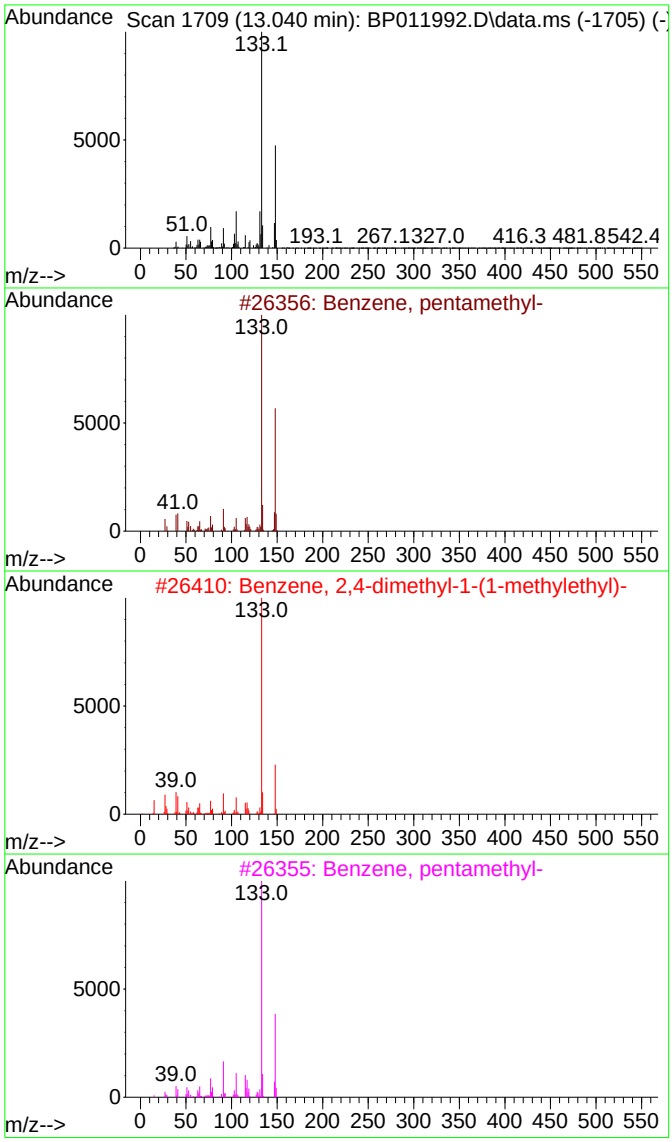
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

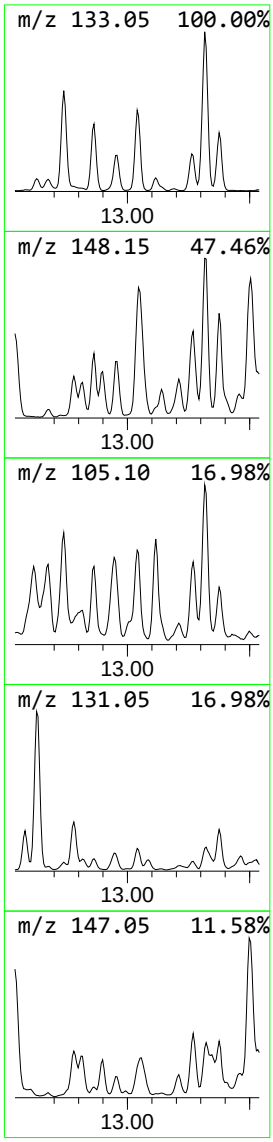
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 Benzene, pentamethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.040	19.74 ng/ul	2200040	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	74
2	Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	72
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	72
4	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	72
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	72





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

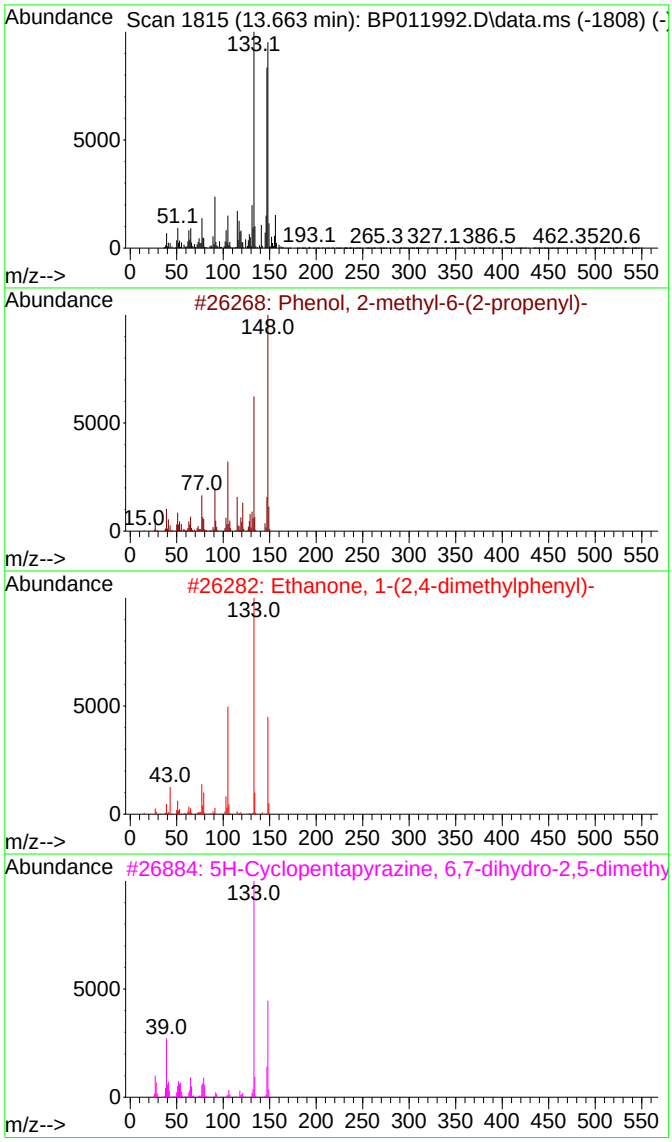
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

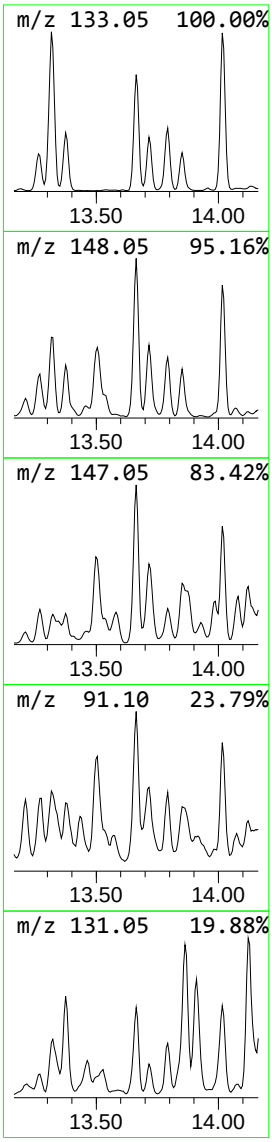
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 35 Phenol, 2-methyl-6-(2-prope... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	50.27 ng/ul	5603650	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methyl-6-(2-propenyl)-	148	C10H12O	003354-58-3	64
2	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	55
3	5H-Cyclopentapyrazine, 6,7-dihyd...	148	C9H12N2	038917-61-2	52
4	1,4-Benzenedicarboxaldehyde, 2-m...	148	C9H8O2	027587-17-3	50
5	Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	49





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

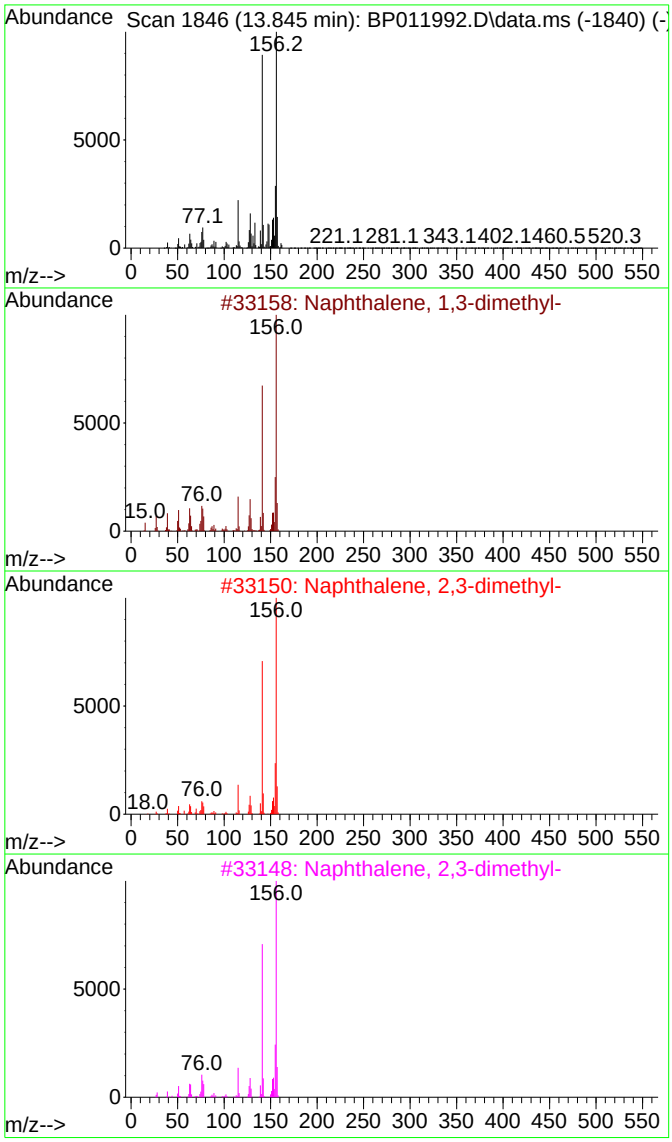
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

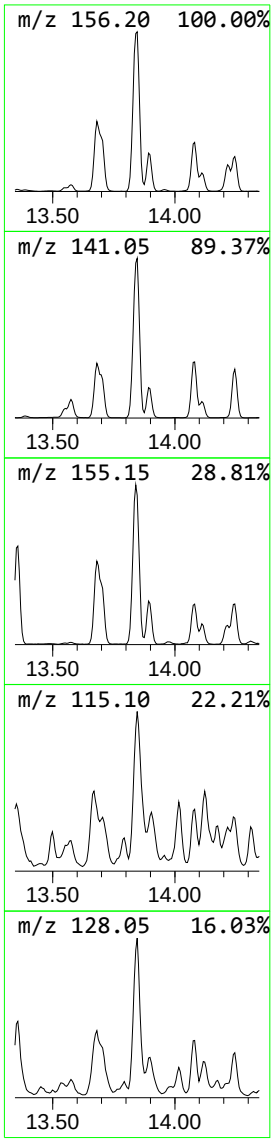
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 36 Naphthalene, 1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.845	62.13 ng/ul	6925180	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

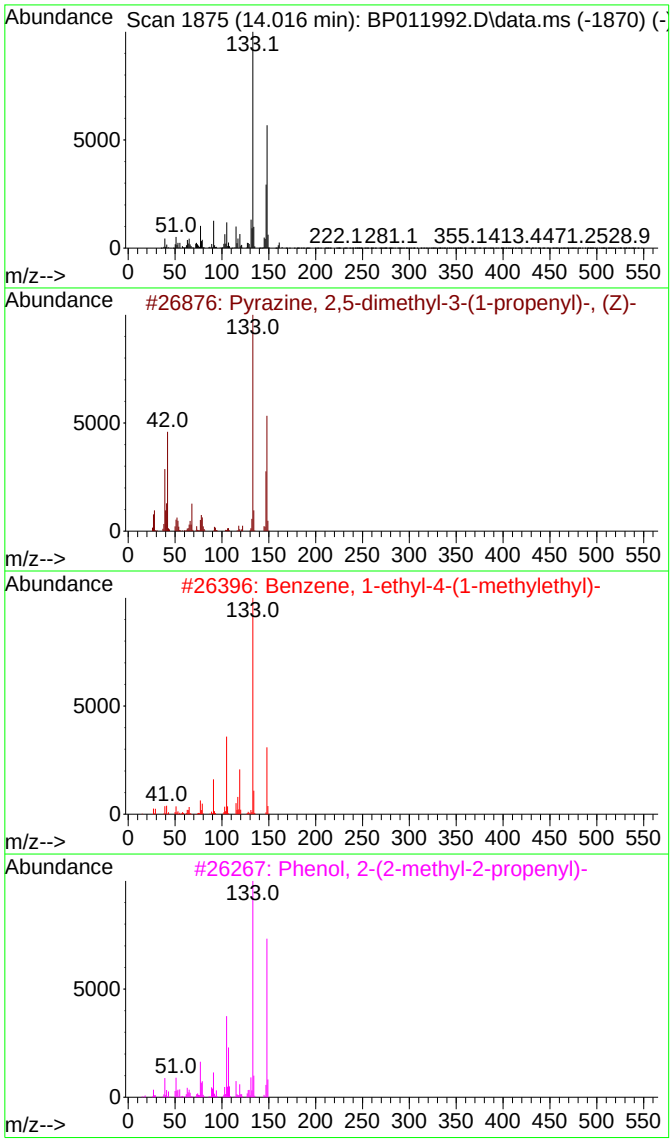
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

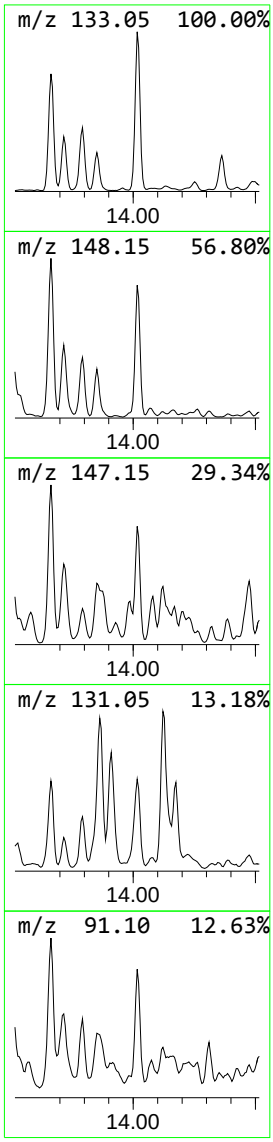
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 Pyrazine, 2,5-dimethyl-3-(1... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.016	27.21 ng/ul	3032600	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrazine, 2,5-dimethyl-3-(1-prop...	148	C9H12N2	055138-73-3	87
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	76
3	Phenol, 2-(2-methyl-2-propenyl)-	148	C10H12O	020944-88-1	72
4	Benzene, pentamethyl-	148	C11H16	000700-12-9	68
5	Benzene, pentamethyl-	148	C11H16	000700-12-9	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
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Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

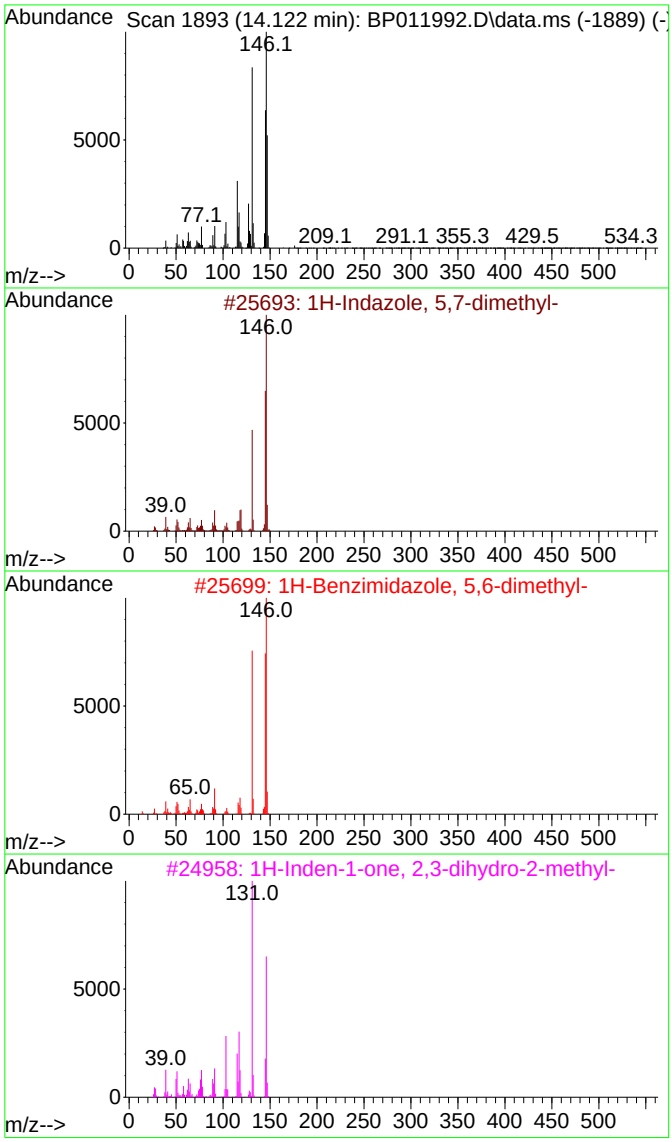
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

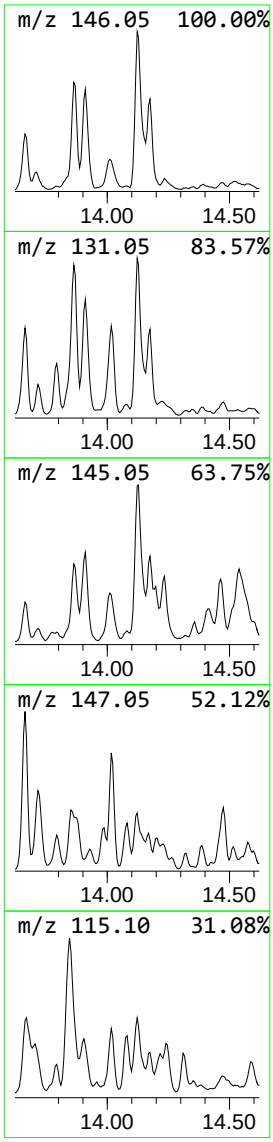
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 1H-Indazole, 5,7-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.122	21.48 ng/ul	2394030	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	76
2	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	58
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	58
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	58
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	58





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

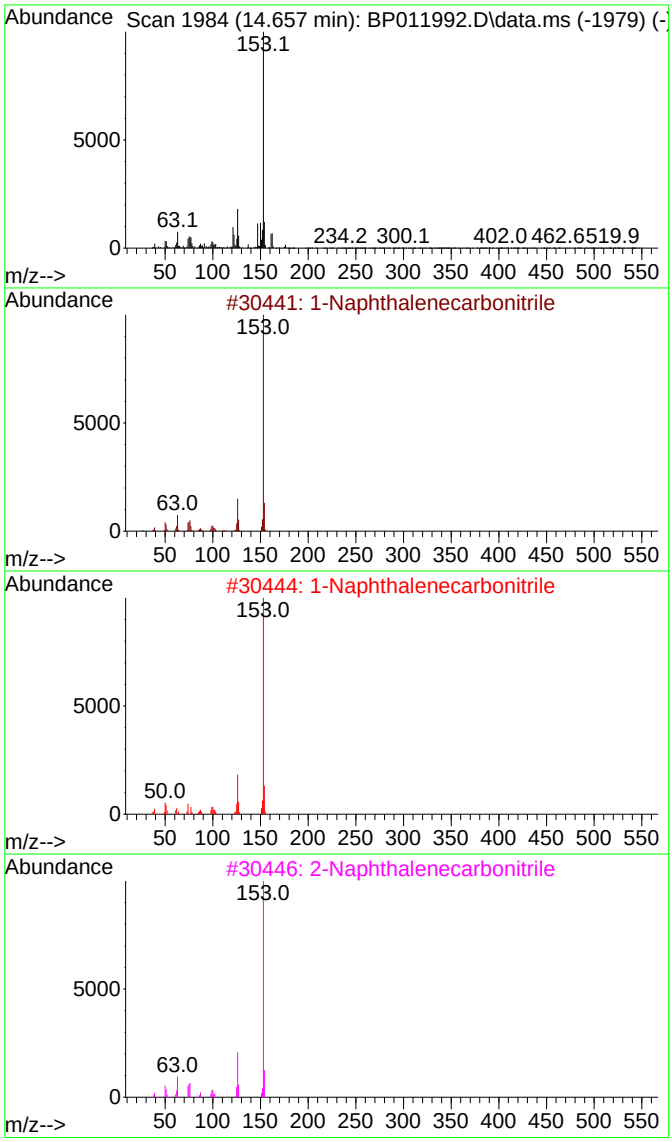
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

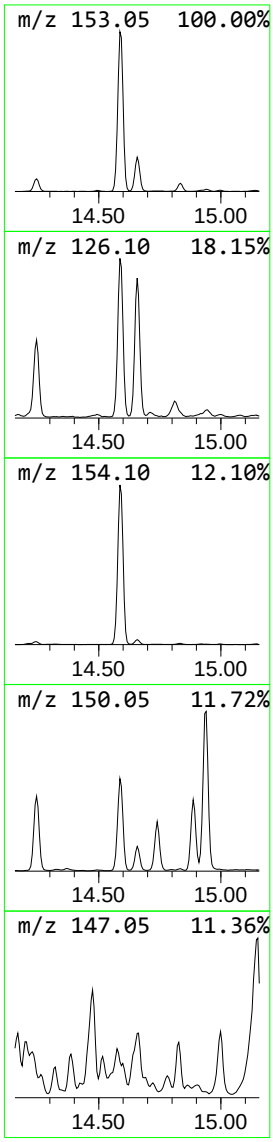
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 1-Naphthalenecarbonitrile Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.657	20.34 ng/ul	2267280	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	95
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	93
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	93
4	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	92
5	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	92





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

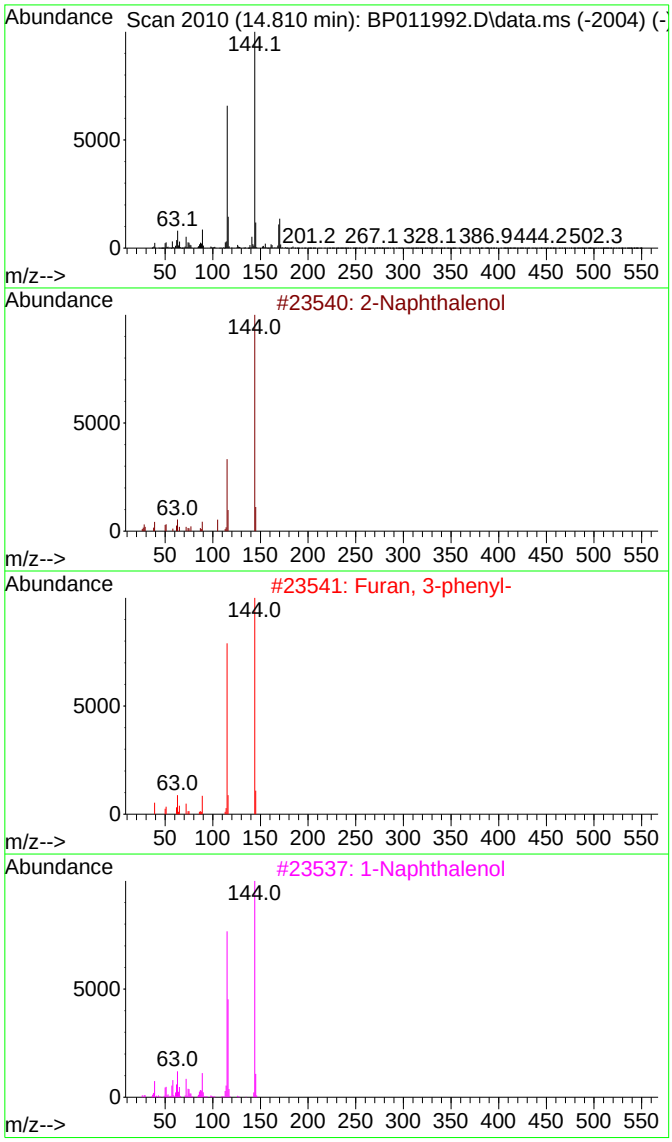
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

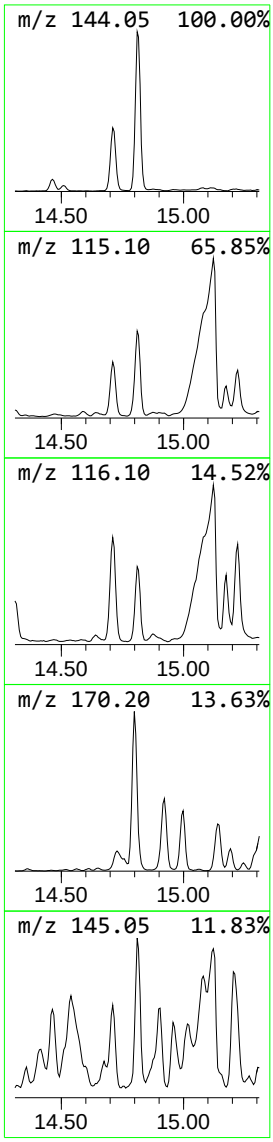
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 40 2-Naphthalenol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.810	37.95 ng/ul	4230090	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenol	144	C10H8O	000135-19-3	87
2	Furan, 3-phenyl-	144	C10H8O	013679-41-9	87
3	1-Naphthalenol	144	C10H8O	000090-15-3	81
4	1-Naphthalenol	144	C10H8O	000090-15-3	81
5	2-Naphthalenol	144	C10H8O	000135-19-3	76





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

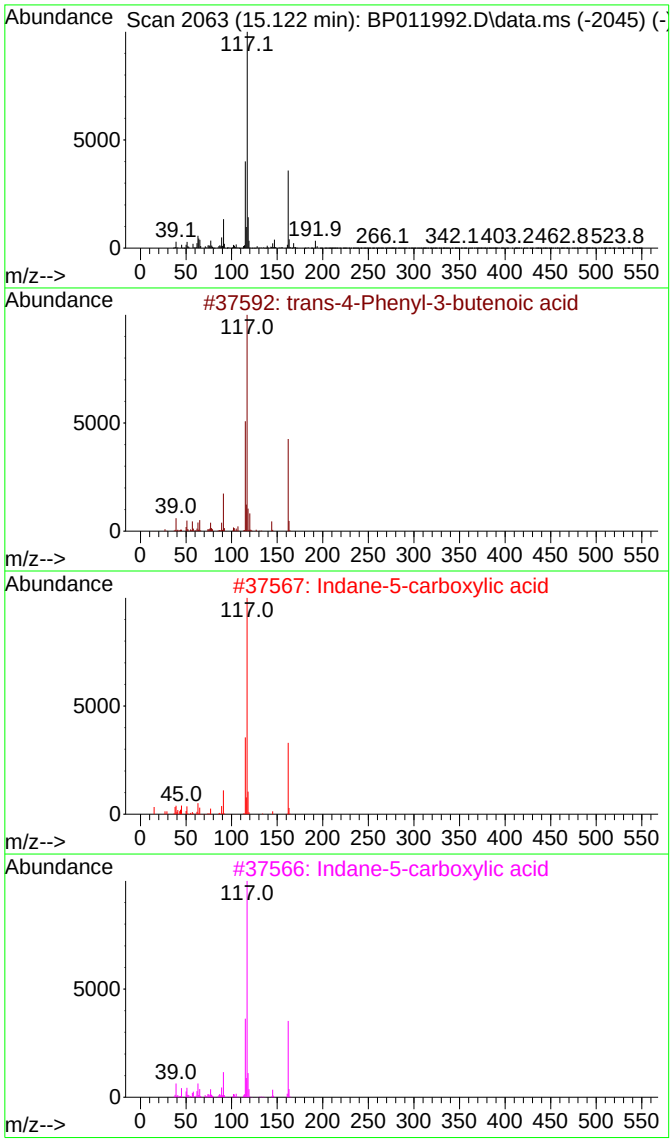
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

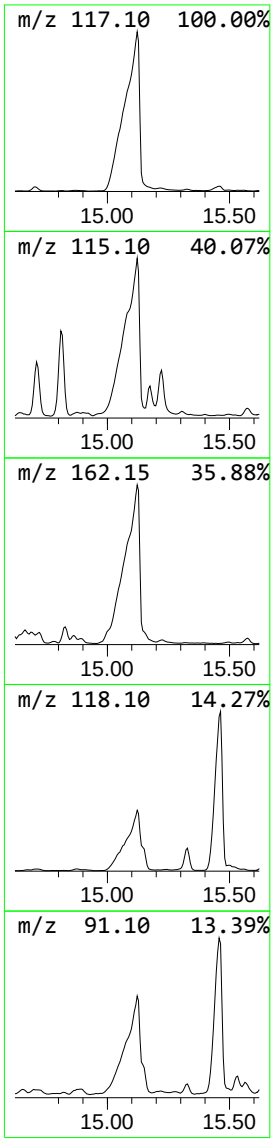
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 trans-4-Phenyl-3-butenoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.122	177.37 ng/ul	19770400	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-4-Phenyl-3-butenoic acid	162	C10H10O2	001914-58-5	94
2	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	91
3	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	91
4	Benzene, 1-(chloromethyl)-3-ethe...	152	C9H9Cl	039833-65-3	64
5	Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

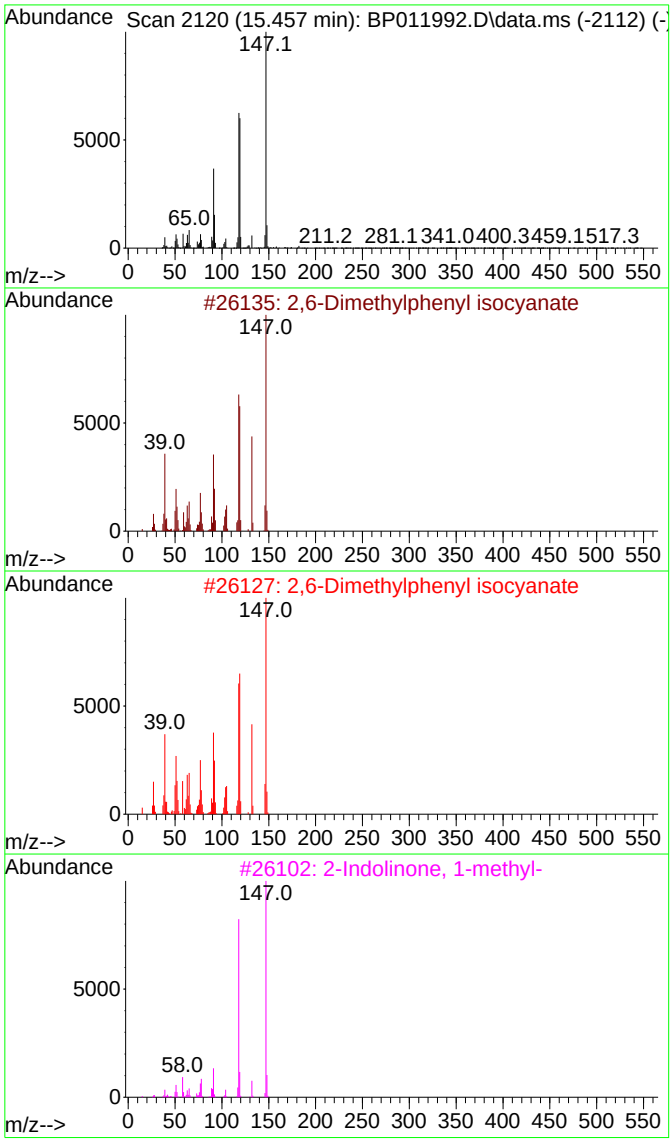
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

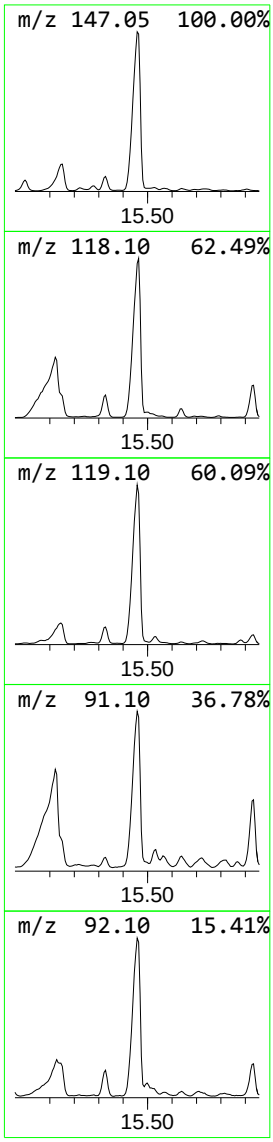
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 42 2,6-Dimethylphenyl isocyanate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.457	95.12 ng/ul	10602300	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	86
2	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	78
3	2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	76
4	2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	68
5	2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	64





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Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

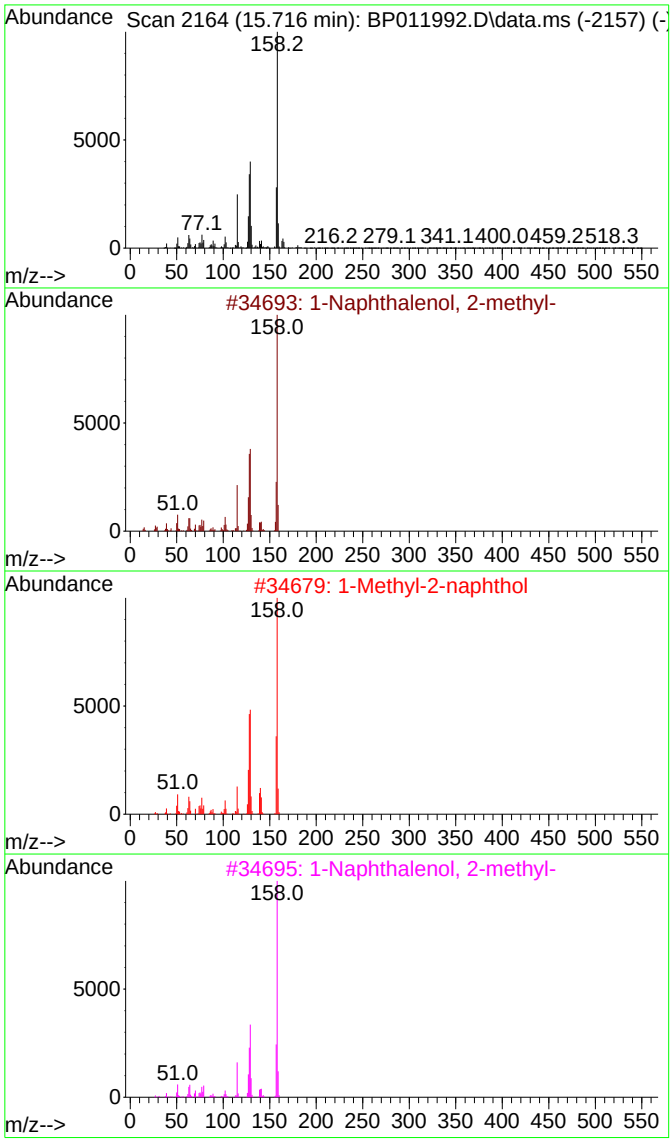
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

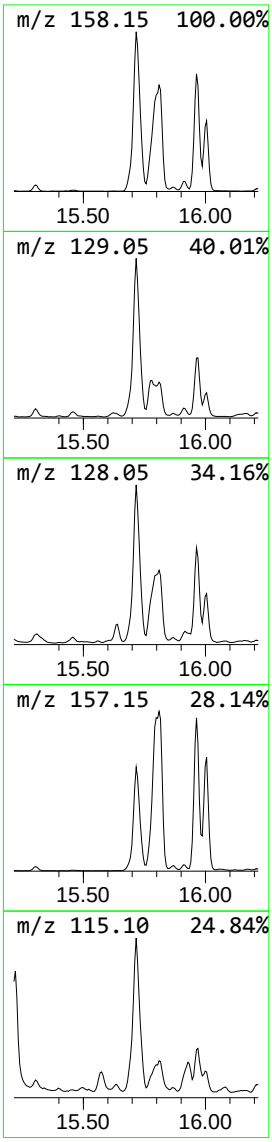
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 43 1-Naphthalenol, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.716	100.38 ng/ul	11189000	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	93
2	1-Methyl-2-naphthol	158	C11H10O	001076-26-2	83
3	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	83
4	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	80
5	7-Methylnaphthalen-2-ol	158	C11H10O	026593-50-0	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

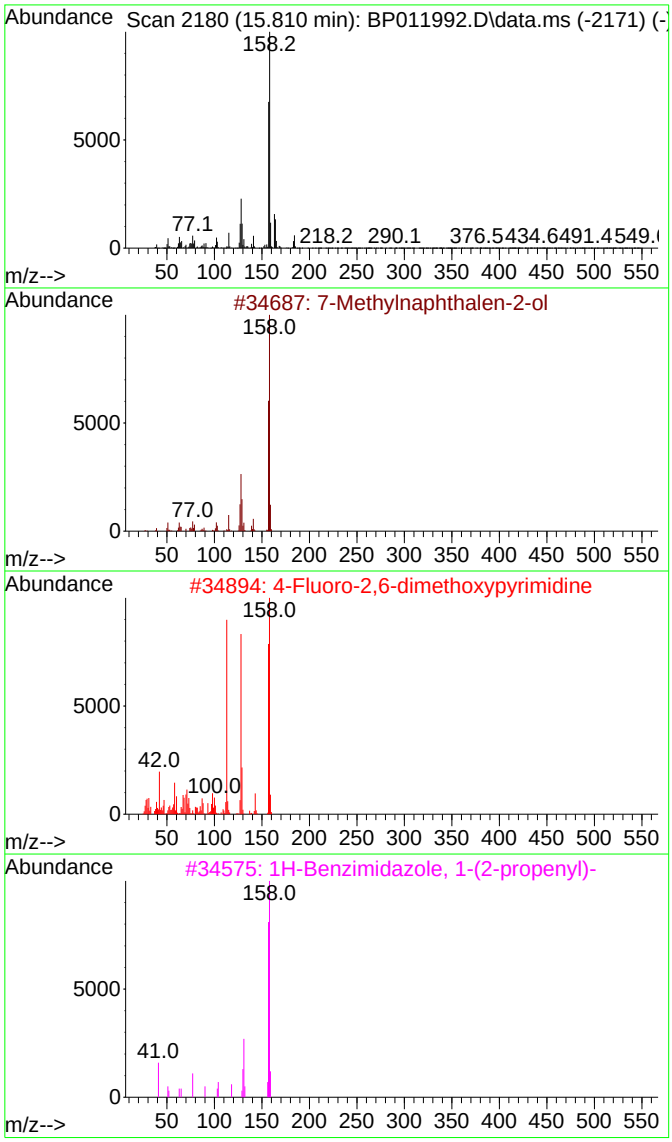
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

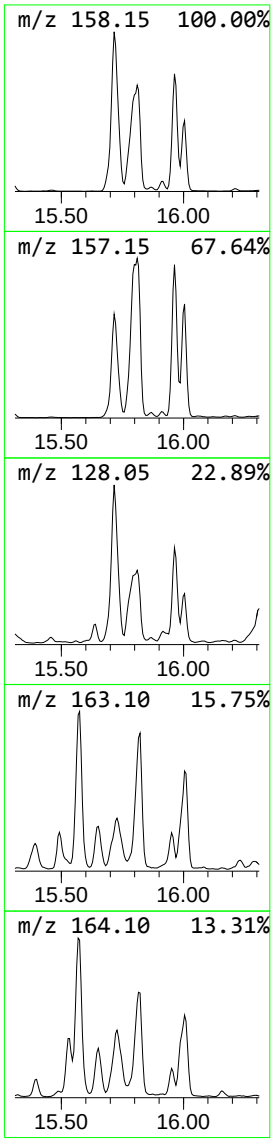
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 44 7-Methylnaphthalen-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.810	82.78 ng/ul	9227220	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methylnaphthalen-2-ol	158	C11H10O	026593-50-0	93
2	4-Fluoro-2,6-dimethoxypyrimidine	158	C6H7FN2O2	000658-87-7	64
3	1H-Benzimidazole, 1-(2-propenyl)-	158	C10H10N2	019018-22-5	50
4	1-(2-Aminophenyl)pyrrole	158	C10H10N2	006025-60-1	50
5	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

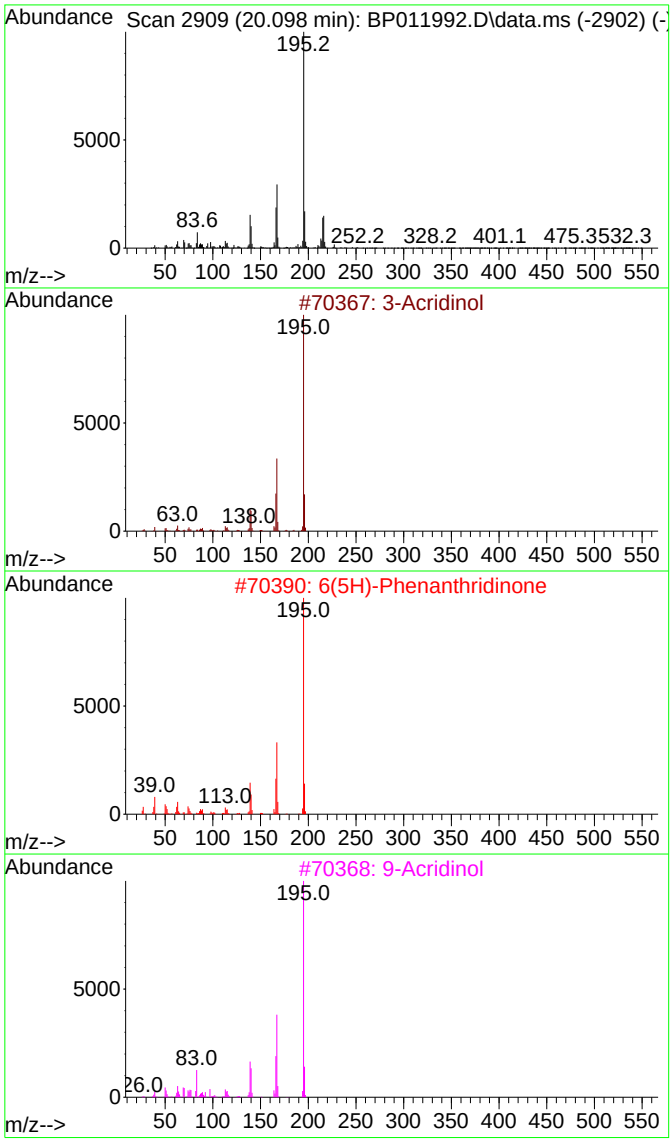
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

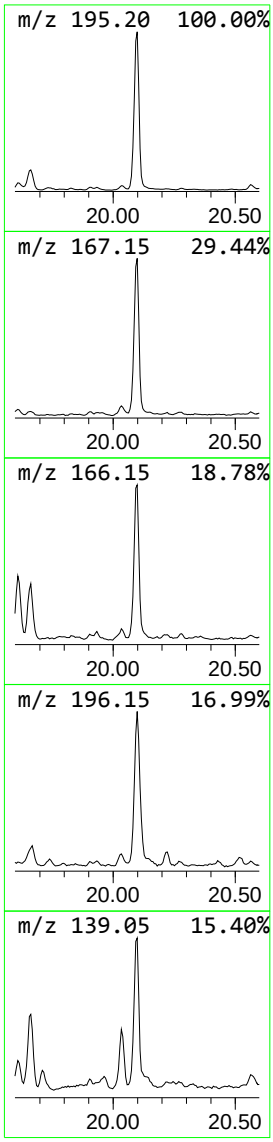
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 61 3-Acridinol Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.098	10.55 ng/ul	3541960	Chrysene-d12	21.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Acridinol	195	C13H9NO	007132-70-9	96
2	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	95
3	9-Acridinol	195	C13H9NO	000643-62-9	95
4	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	94
5	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011992.D
Acq On : 07 Oct 2022 04:19
Operator : CG/JU
Sample : N4923-09
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E10011

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.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
Butane, 2-metho...	3.034	54.5	ng/ul	5828360	1	7.875	2140350	20.0
Benzene, 1,2,3-...	7.516	21.4	ng/ul	2291290	1	7.875	2140350	20.0
Benzofuran	7.593	37.3	ng/ul	3994050	1	7.875	2140350	20.0
Indane	8.252	226.0	ng/ul	24184300	1	7.875	2140350	20.0
Indene	8.416	120.6	ng/ul	12909000	1	7.875	2140350	20.0
p-Aminotoluene	8.781	27.8	ng/ul	2976050	1	7.875	2140350	20.0
Benzenamine, 3-...	8.869	41.0	ng/ul	4382870	1	7.875	2140350	20.0
Benzofuran, 2-m...	9.234	21.7	ng/ul	2323540	1	7.875	2140350	20.0
Phenol, 2,6-dim...	9.316	23.7	ng/ul	21033900	2	10.687	17778400	20.0
Phenol, 3,5-dim...	10.181	22.1	ng/ul	19654100	2	10.687	17778400	20.0
Phenol, 3-ethyl-	10.563	8.1	ng/ul	7200180	2	10.687	17778400	20.0
Phenol, 2,3,6-t...	11.281	12.5	ng/ul	11089300	2	10.687	17778400	20.0
Phenol, 3-ethyl...	11.534	14.9	ng/ul	13255500	2	10.687	17778400	20.0
Phenol, 3,4,5-t...	11.722	15.4	ng/ul	13652400	2	10.687	17778400	20.0
Phenol, 2,4,6-t...	11.775	10.5	ng/ul	9330900	2	10.687	17778400	20.0
Phenol, 3,5-die...	12.675	25.5	ng/ul	2842850	3	14.522	2229350	20.0
unknown-01	12.722	26.0	ng/ul	2902660	3	14.522	2229350	20.0
Ethanone, 1-(2-...	12.945	23.5	ng/ul	2619520	3	14.522	2229350	20.0
Benzene, pentam...	13.040	19.7	ng/ul	2200040	3	14.522	2229350	20.0
Phenol, 2-methy...	13.663	50.3	ng/ul	5603650	3	14.522	2229350	20.0
Naphthalene, 1,...	13.845	62.1	ng/ul	6925180	3	14.522	2229350	20.0
Pyrazine, 2,5-d...	14.016	27.2	ng/ul	3032600	3	14.522	2229350	20.0
1H-Indazole, 5,...	14.122	21.5	ng/ul	2394030	3	14.522	2229350	20.0
1-Naphthaleneca...	14.657	20.3	ng/ul	2267280	3	14.522	2229350	20.0
2-Naphthalenol	14.810	38.0	ng/ul	4230090	3	14.522	2229350	20.0
trans-4-Phenyl-...	15.122	177.4	ng/ul	19770400	3	14.522	2229350	20.0
2,6-Dimethylphe...	15.457	95.1	ng/ul	10602300	3	14.522	2229350	20.0
1-Naphthalenol,...	15.716	100.4	ng/ul	11189000	3	14.522	2229350	20.0
7-Methylnaphtha...	15.810	82.8	ng/ul	9227220	3	14.522	2229350	20.0
3-Acridinol	20.098	10.6	ng/ul	3541960	5	21.369	6717540	20.0