LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014500.D
Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB996

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-BP032823.M
Title : SVOA CALIBRATION

Signal : TIC: BP014500.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.381	30	33	36	rBV	21167	26823	0.81%	0.067%
2	3.417	36	39	47	rVB	34928	53129	1.61%	0.133%
3	3.822	105	108	120	rBV	167745	289569	8.79%	0.723%
4	4.040	140	145	147	rBV3	8575	12242	0.37%	0.031%
5	4.140	158	162	168	rVB	8801	13339	0.40%	0.033%
6	4.252	177	181	182	rBV4	5920	8380	0.25%	0.021%
7	4.522	222	227	231	rVV7	4611	8977	0.27%	0.022%
8	4.599	235	240	246	rVB	30418	50616	1.54%	0.126%
9	5.069	316	320	326	rBV2	9856	15510	0.47%	0.039%
10	5.281	354	356	361	rBV6	5117	8209	0.25%	0.020%
11	5.746	430	435	439	rBV7	9627	19307	0.59%	0.048%
12	6.469	555	558	562	rBV5	5055	8131	0.25%	0.020%
13	6.634	581	586	588	rVV6	5469	9076	0.28%	0.023%
14	7.181	671	679	689	rBV	75916	168796	5.12%	0.421%
15	7.334	695	705	716	rBV	328285	552651	16.77%	1.379%
16	7.487	725	731	734	rVV4	26950	45609	1.38%	0.114%
17	7.540	734	740	754	rVV	321007	568483	17.25%	1.419%
18	7.658	755	760	768	rVV9	8584	21717	0.66%	0.054%
19	7.899	798	801	808	rVB9	5259	9980	0.30%	0.025%
20	8.005	813	819	827	rVV	503413	813525	24.69%	2.031%
21	8.093	829	834	843	rVV8	16326	43600	1.32%	0.109%
22	8.305	866	870	877	rBV9	5053	11041	0.34%	0.028%
23	8.381	877	883	890	rVB	17705	31032	0.94%	0.077%
24	8.493	897	902	905	rVB7	4759	8166	0.25%	0.020%
25	8.546	907	911	917	rVB9	4783	9052	0.27%	0.023%
26	8.669	925	932	937	rBV	15798	30853	0.94%	0.077%
27	8.734	937	943	952	rBV	274685	495972	15.05%	1.238%
28	9.010	984	990	999	rBV2	61966	115659	3.51%	0.289%
29	9.122	1004	1009	1013	rVV5	15009	27889	0.85%	0.070%
30	9.181	1013	1019	1026	rVV	430682	743684	22.57%	1.856%
31	9.252	1026	1031	1039	rVB	115662	187865	5.70%	0.469%
32	9.446	1060	1064	1070	rBV9	5786	12334	0.37%	0.031%
33	9.534	1072	1079	1087	rBV5	23597	60518	1.84%	0.151%
34	9.646	1093	1098	1102	rBV4	11397	20668	0.63%	0.052%
35	9.757	1112	1117	1123	rVB6	14762	27007	0.82%	0.067%
36	9.828	1123	1129	1136	rBV6	5757	13312	0.40%	0.033%

LSC Area Percent Report

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Operator : CG/JU
Sample : 02093-14
Misc :
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Instrument :
BNA_P
ClientSampleId :
CB996

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-BP032823.M
Title : SVOA CALIBRATION

37	9.910	1136	1143	1154	rBV	358238	660719	20.05%	1.649%
38	10.034	1157	1164	1169	rBV7	16520	37675	1.14%	0.094%
39	10.093	1171	1174	1176	rBV4	7538	9249	0.28%	0.023%
40	10.234	1187	1198	1205	rBV4	53794	111481	3.38%	0.278%
41	10.304	1207	1210	1213	rVV4	5854	9922	0.30%	0.025%
42	10.363	1213	1220	1229	rVB4	41195	107221	3.25%	0.268%
43	10.457	1229	1236	1254	rBV	480489	1019423	30.93%	2.544%
44	10.610	1257	1262	1267	rVB8	9418	20948	0.64%	0.052%
45	10.828	1287	1299	1306	rVV	755880	1321396	40.10%	3.298%
46	10.875	1306	1307	1313	rVV2	26027	33214	1.01%	0.083%
47	10.981	1314	1325	1338	rVV	452026	989070	30.01%	2.469%
48	11.081	1338	1342	1353	rVV6	22609	58028	1.76%	0.145%
49	11.181	1355	1359	1364	rVB8	7701	14617	0.44%	0.036%
50	11.422	1391	1400	1405	rBV8	12436	30339	0.92%	0.076%
51	11.551	1419	1422	1426	rBV6	7130	10846	0.33%	0.027%
52	11.651	1432	1439	1442	rBV9	16561	35850	1.09%	0.089%
53	11.816	1464	1467	1470	rVB5	6367	7298	0.22%	0.018%
54	12.028	1499	1503	1510	rVB6	16071	28455	0.86%	0.071%
55	12.081	1510	1512	1519	rVB8	5840	8222	0.25%	0.021%
56	12.175	1519	1528	1533	rBV	72030	134033	4.07%	0.335%
57	12.222	1533	1536	1543	rVB2	27991	44861	1.36%	0.112%
58	12.351	1550	1558	1564	rBV9	19163	66853	2.03%	0.167%
59	12.540	1587	1590	1596	rVB7	14964	24956	0.76%	0.062%
60	12.704	1615	1618	1624	rVB5	18939	28343	0.86%	0.071%
61	12.822	1633	1638	1641	rBV2	27121	42579	1.29%	0.106%
62	12.928	1650	1656	1660	rBV9	15985	29272	0.89%	0.073%
63	13.010	1667	1670	1671	rBV3	8608	7962	0.24%	0.020%
64	13.257	1709	1712	1720	rVB10	12912	25710	0.78%	0.064%
65	13.363	1726	1730	1732	rBV5	6827	11364	0.34%	0.028%
66	13.457	1743	1746	1751	rBV3	25540	32979	1.00%	0.082%
67	13.540	1757	1760	1764	rVB6	24003	35236	1.07%	0.088%
68	13.710	1785	1789	1798	rVB8	23127	43038	1.31%	0.107%
69	14.069	1841	1850	1857	rBV	1237708	1788326	54.27%	4.464%
70	14.222	1872	1876	1878	rBV4	16181	24736	0.75%	0.062%
71	14.357	1892	1899	1907	rBV	1253908	1918719	58.22%	4.789%
72	14.475	1914	1919	1925	rBV4	29455	59800	1.81%	0.149%
73	14.581	1932	1937	1944	rBV2	80805	126868	3.85%	0.317%
74	14.657	1944	1950	1957	rVV	1218430	1859796	56.44%	4.642%
75	14.722	1957	1961	1972	rVB	75506	124544	3.78%	0.311%
76	14.934	1992	1997	2001	rBV4	42834	88043	2.67%	0.220%

LSC Area Percent Report

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Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Title : SVOA CALIBRATION

77	15.310	2057	2061	2065	rBV	42835	57146	1.73%	0.143%
78	15.398	2071	2076	2085	rBV	419407	566737	17.20%	1.415%
79	15.481	2088	2090	2095	rVB3	28605	36534	1.11%	0.091%
80	15.657	2114	2120	2126	rBV	1694754	2438386	73.99%	6.086%
81	15.710	2126	2129	2135	rVB	39160	57429	1.74%	0.143%
82	15.787	2137	2142	2153	rBV	371636	655060	19.88%	1.635%
83	15.869	2153	2156	2159	rVV	34840	38453	1.17%	0.096%
84	16.022	2177	2182	2193	rVB	261528	390803	11.86%	0.975%
85	16.181	2203	2209	2219	rVB	490649	705004	21.39%	1.760%
86	16.510	2263	2265	2269	rVB3	23077	26075	0.79%	0.065%
87	16.592	2276	2279	2285	rVB4	35513	59464	1.80%	0.148%
88	16.692	2291	2296	2300	rBV	264704	376075	11.41%	0.939%
89	17.151	2371	2374	2377	rBV2	24163	27092	0.82%	0.068%
90	17.422	2415	2420	2431	rVB	1662114	2336237	70.89%	5.831%
91	17.522	2431	2437	2445	rBV2	1933545	2770221	84.06%	6.914%
92	18.286	2563	2567	2582	rBV	705665	1117304	33.90%	2.789%
93	19.051	2694	2697	2701	rVB3	66437	77421	2.35%	0.193%
94	19.516	2773	2776	2783	rBV	149442	217859	6.61%	0.544%
95	19.751	2811	2816	2821	rBV	2595742	3295460	100.00%	8.225%
96	19.792	2821	2823	2830	rVB	107594	134269	4.07%	0.335%
97	21.186	3056	3060	3067	rBV	786976	1055207	32.02%	2.634%
98	21.510	3110	3115	3122	rBV	1900279	2561652	77.73%	6.394%
99	23.863	3508	3515	3524	rVB2	1587357	3190813	96.82%	7.964%
100	24.027	3537	3543	3554	rVB2	1104542	2298900	69.76%	5.738%

Sum of corrected areas: 40064313

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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Misc :
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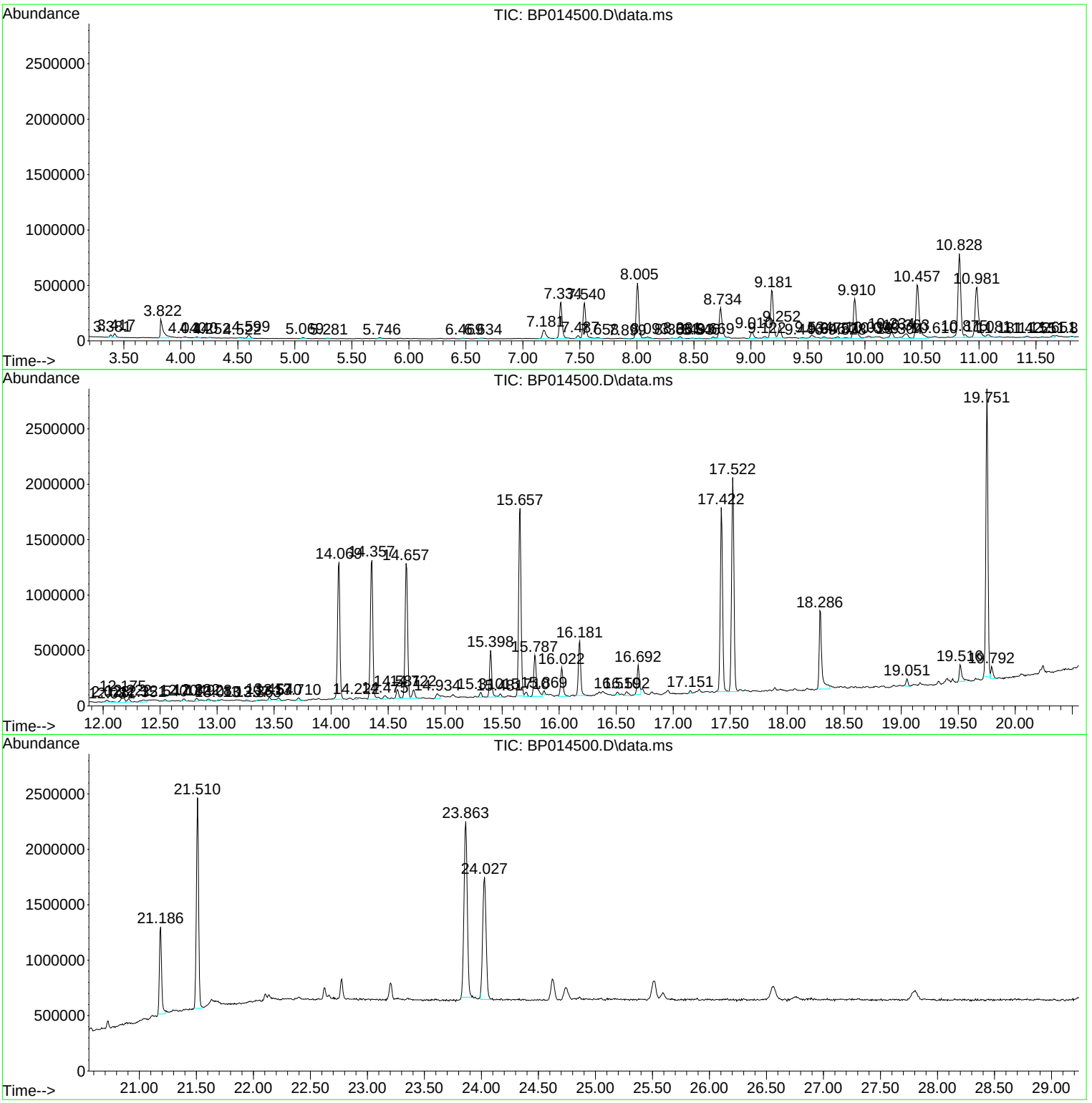
Instrument :
BNA_P
ClientSampleId :
CB996

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014500.D
Acq On : 03 Apr 2023 22:36
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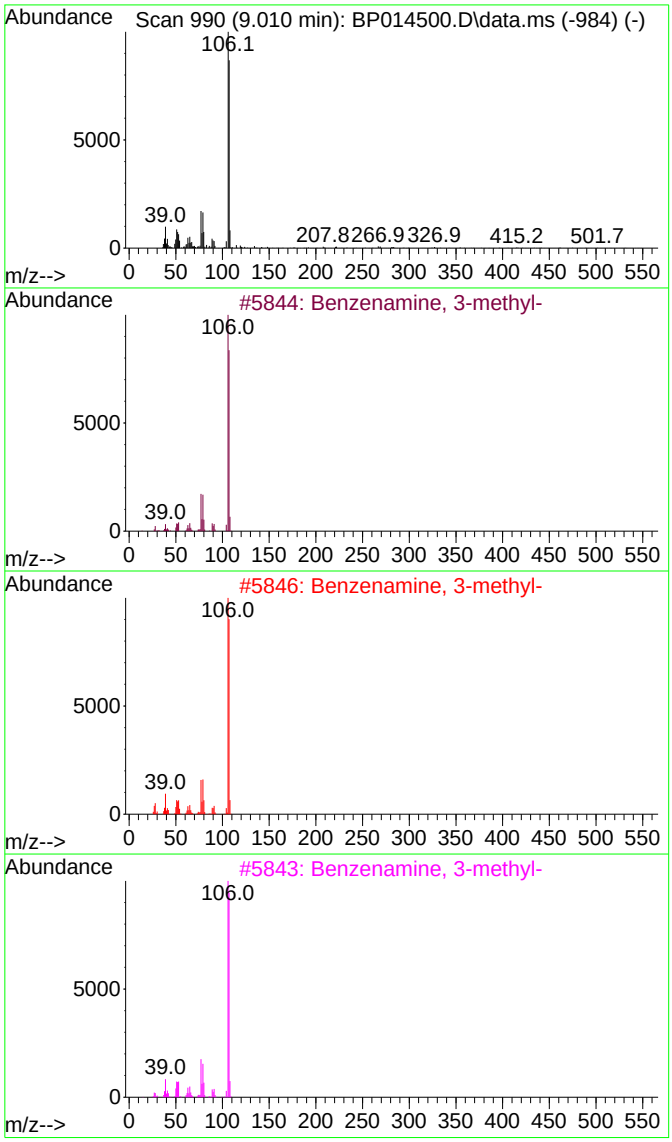
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

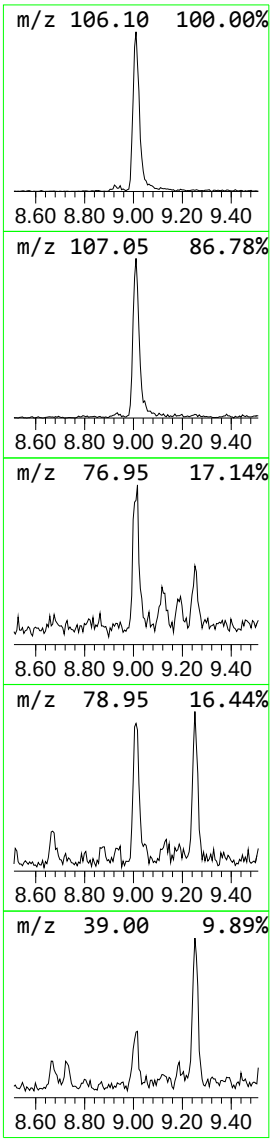
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzenamine, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.010	2.84 ng/ul	115659	1,4-Dichlorobenzene-d4	8.005

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	96
5	o-Toluidine	107	C7H9N	000095-53-4	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014500.D
Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

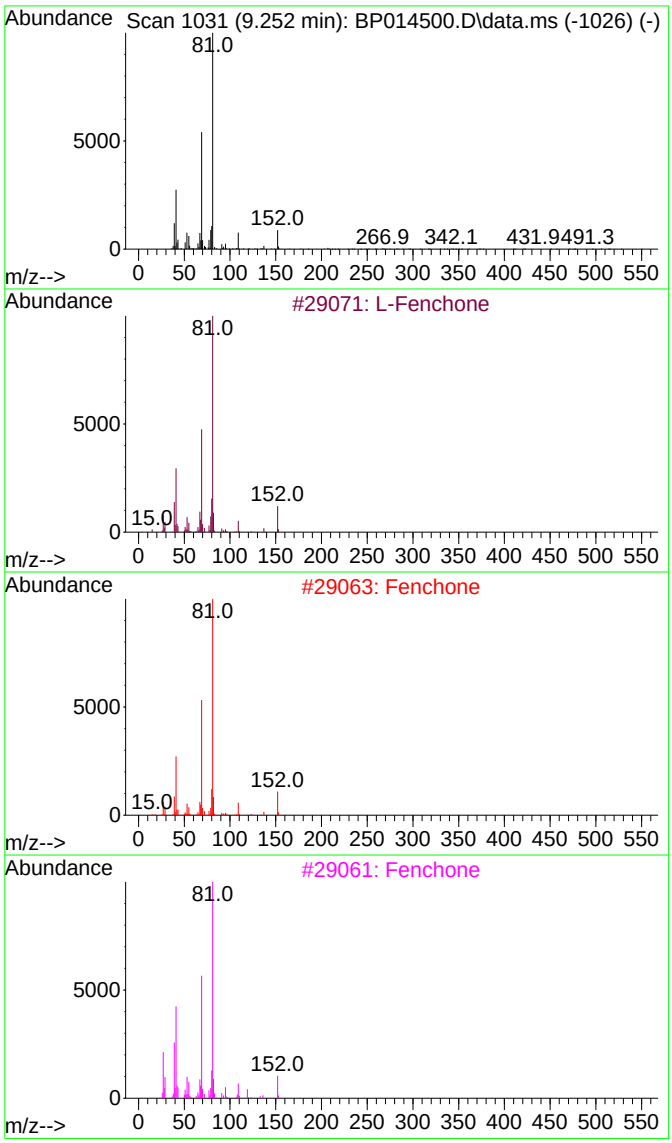
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

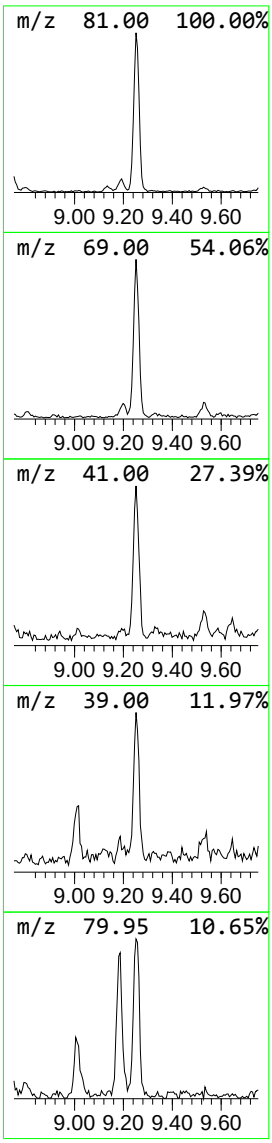
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 L-Fenchone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.252	4.62 ng/ul	187865	1,4-Dichlorobenzene-d4	8.005

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	L-Fenchone	152	C10H16O	007787-20-4	90
2	Fenchone	152	C10H16O	001195-79-5	90
3	Fenchone	152	C10H16O	001195-79-5	78
4	Benzene, 1-[(2-bromocyclohexyl)t...	360	C12H13BrN2O4S	1000362-21-3	42
5	1-[4-(4-tert-Butyl-spirocyclohex...	371	C20H28F3NO2	1000301-43-3	36





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CB996

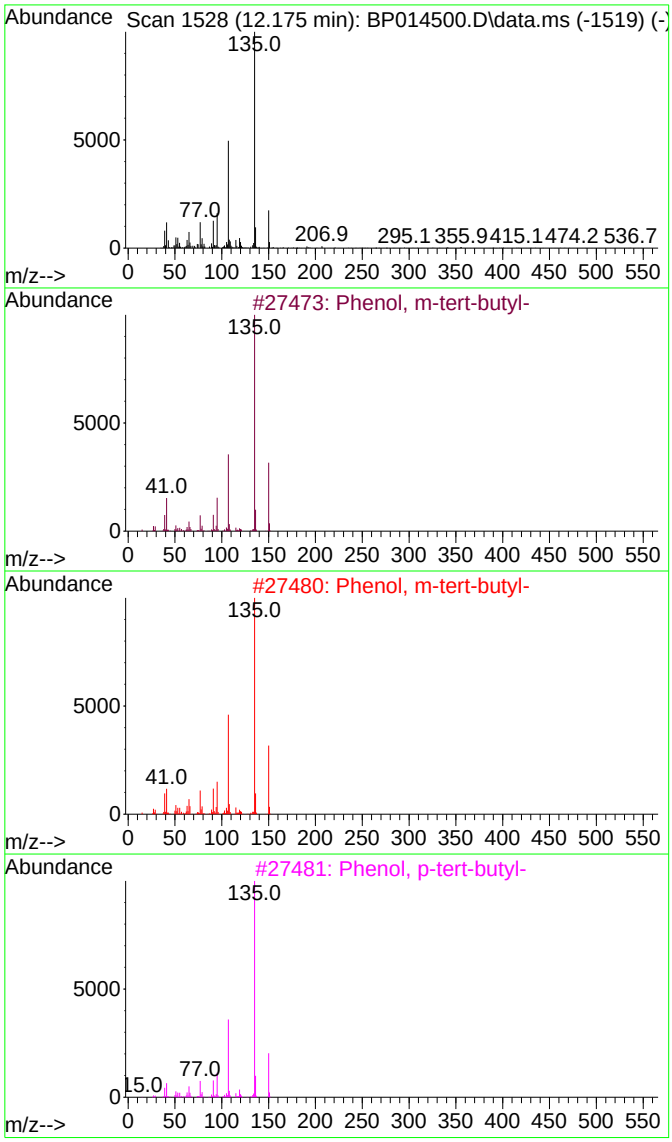
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

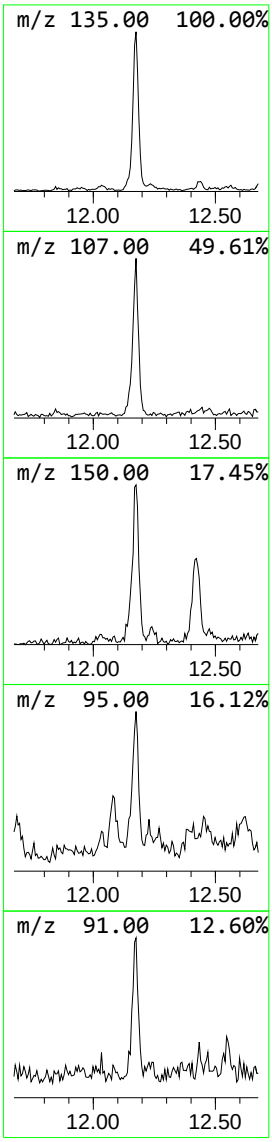
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, m-tert-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.175	2.03 ng/ul	134033	Naphthalene-d8	10.828

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





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

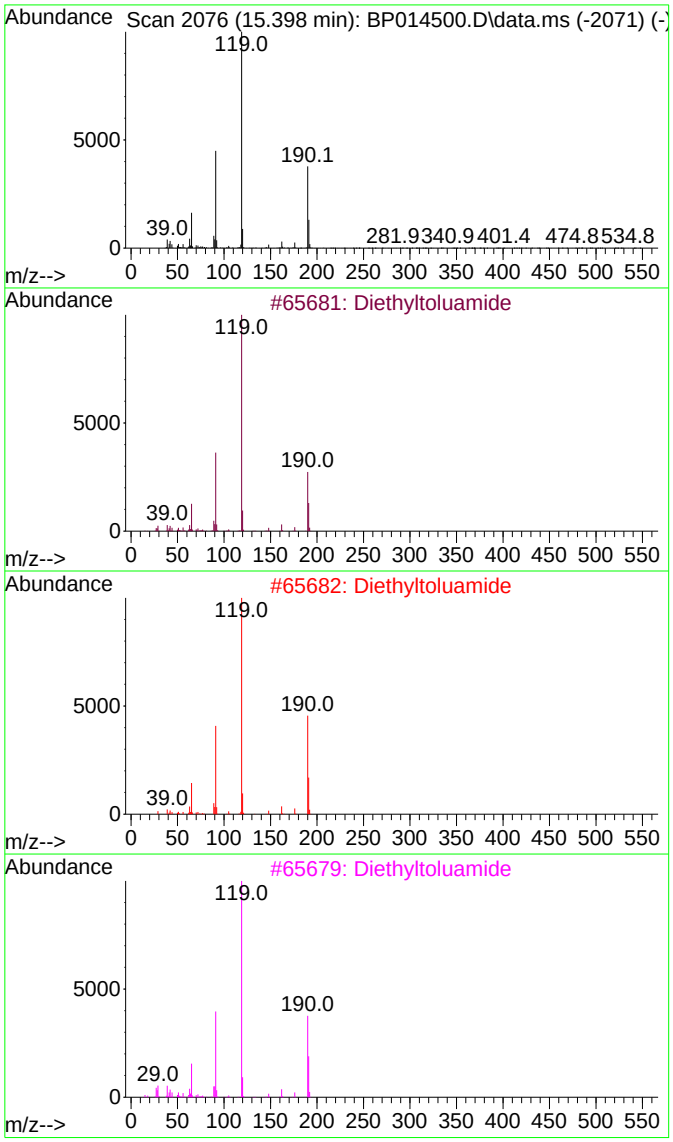
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

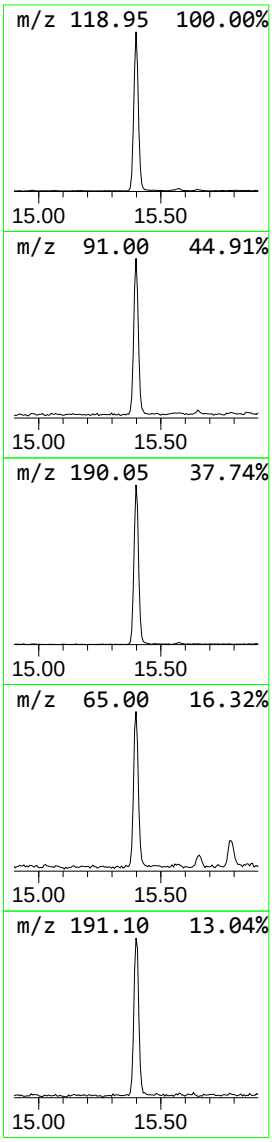
****

Peak Number 4 Diethyltoluamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.398	6.09 ng/ul	566737	Acenaphthene-d10	14.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	96
3	Diethyltoluamide	191	C12H17NO	000134-62-3	95
4	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	90
5	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	87





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Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

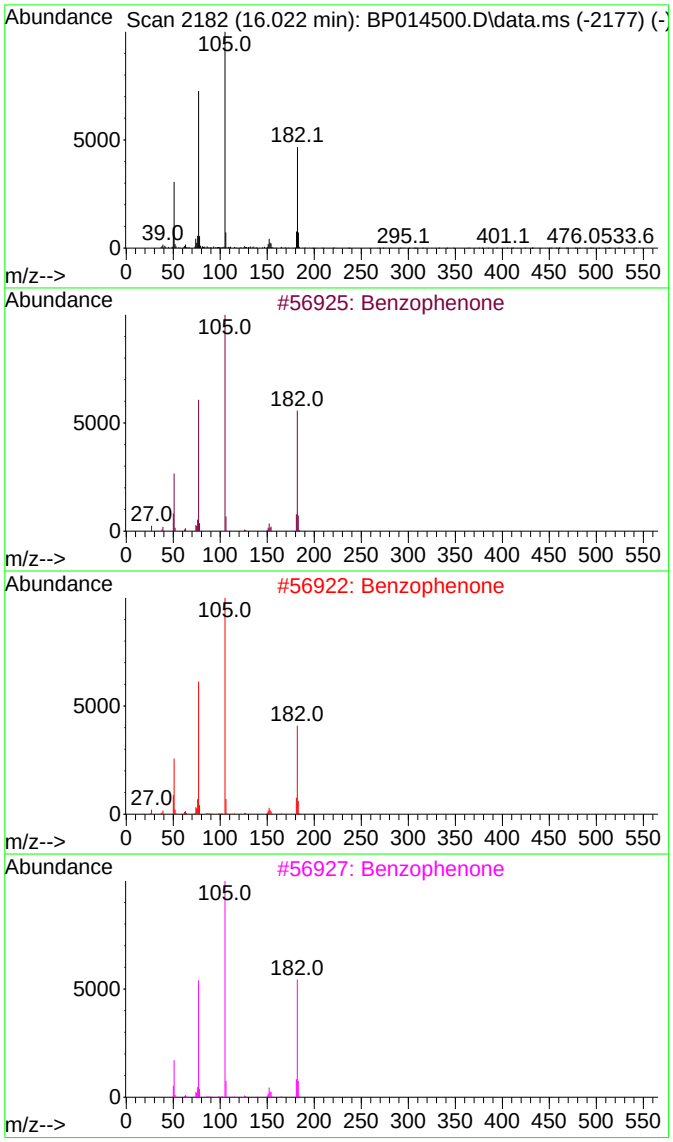
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

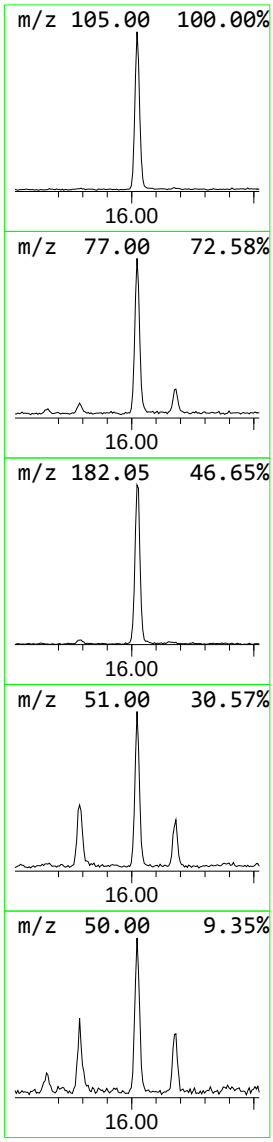
****

Peak Number 5 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.022	4.20 ng/ul	390803	Acenaphthene-d10	14.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	91
5	1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78





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Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

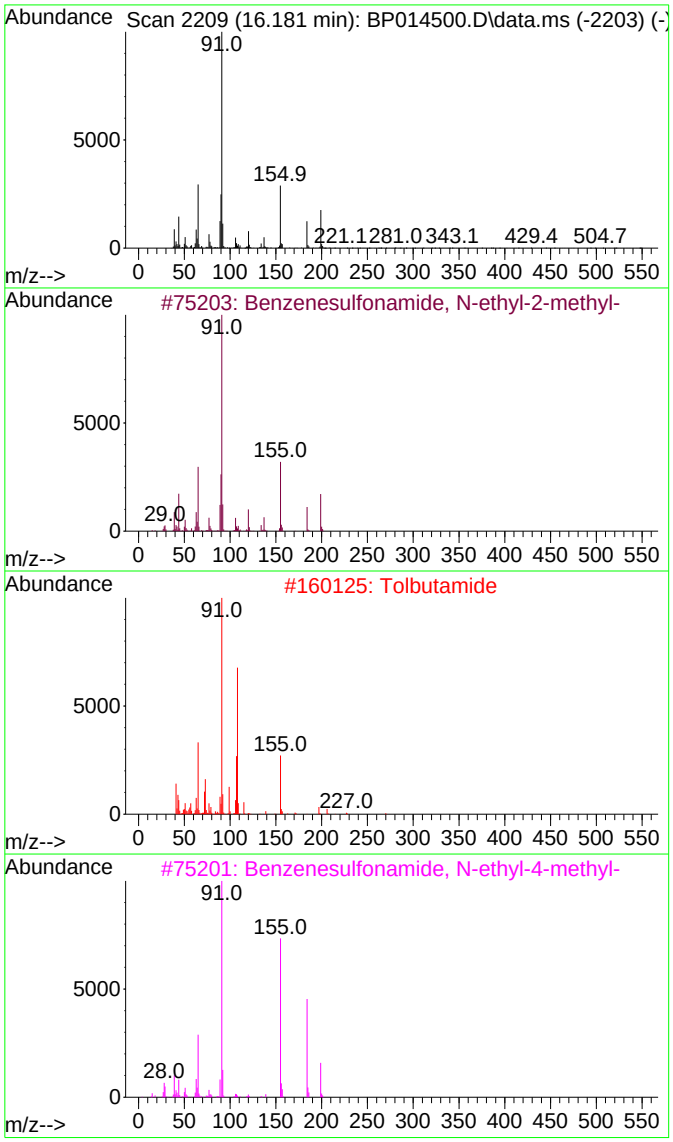
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

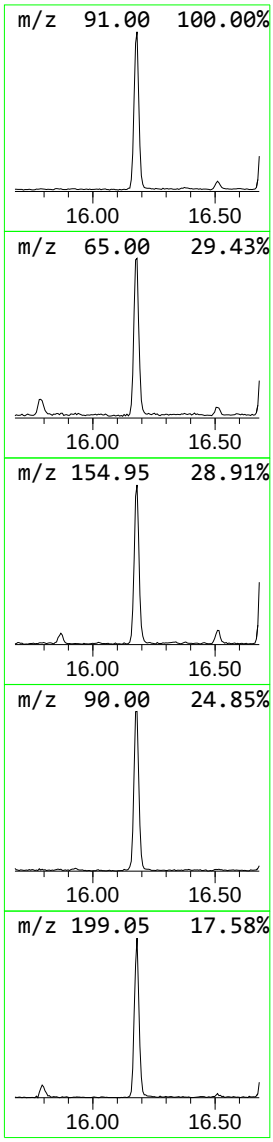
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.181	6.04 ng/ul	705004	Phenanthrene-d10	17.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	49
4	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47
5	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	47





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Data File : BP014500.D
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Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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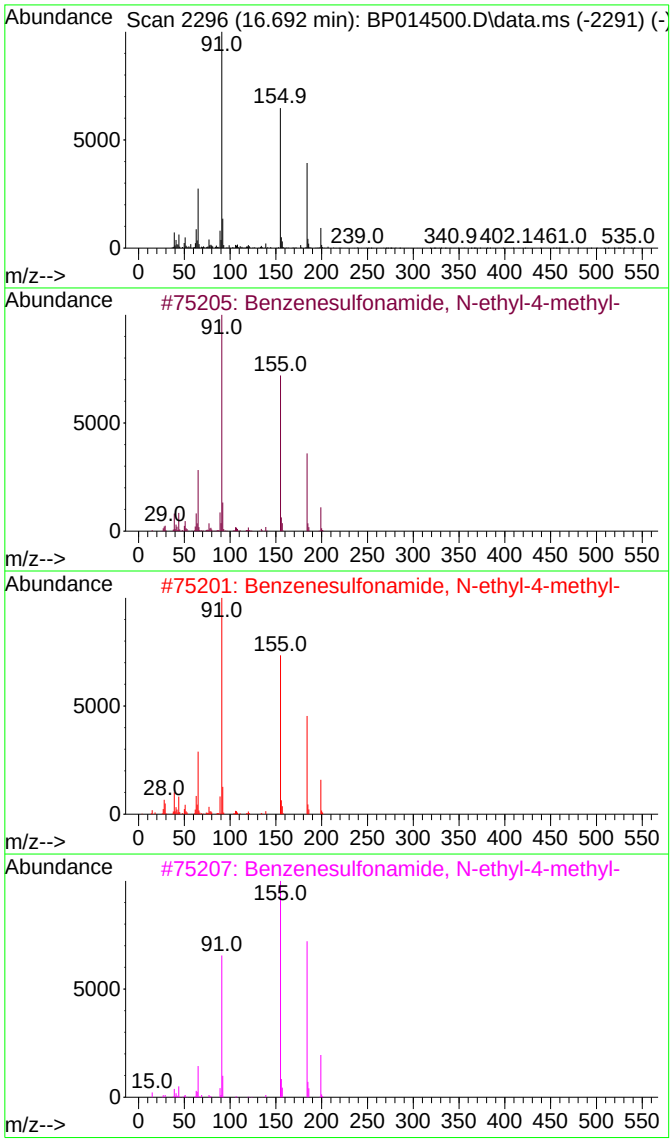
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

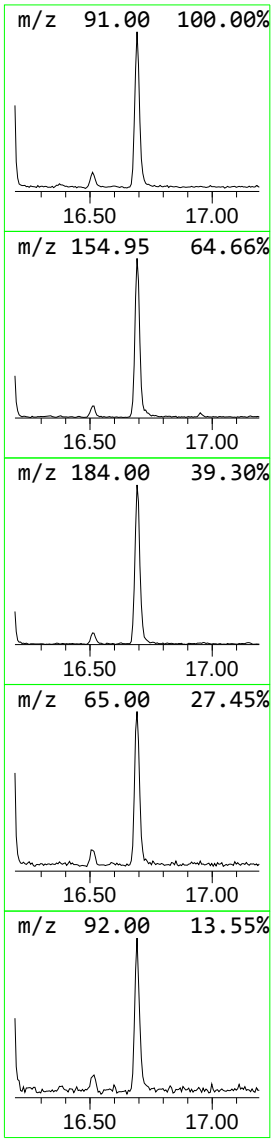
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.692	3.22 ng/ul	376075	Phenanthrene-d10	17.422

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	96
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4	N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	86
5	Benzene, 4-(ethylsulfonyl)-1-met...	184	C9H12O2S	007569-34-8	80





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Data File : BP014500.D
Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

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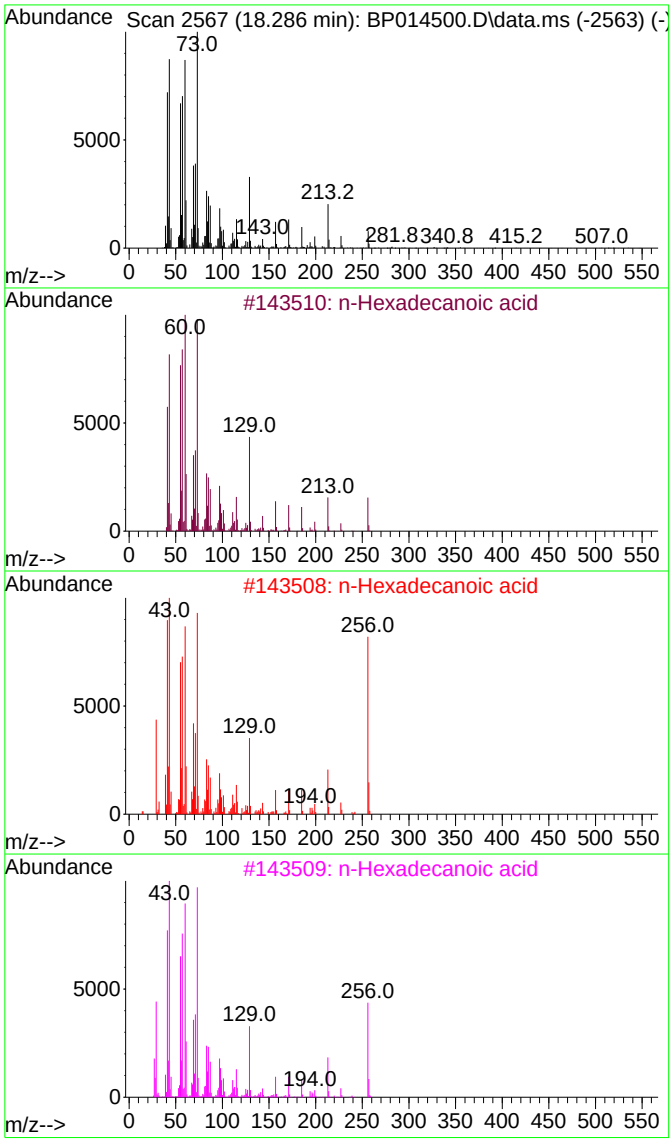
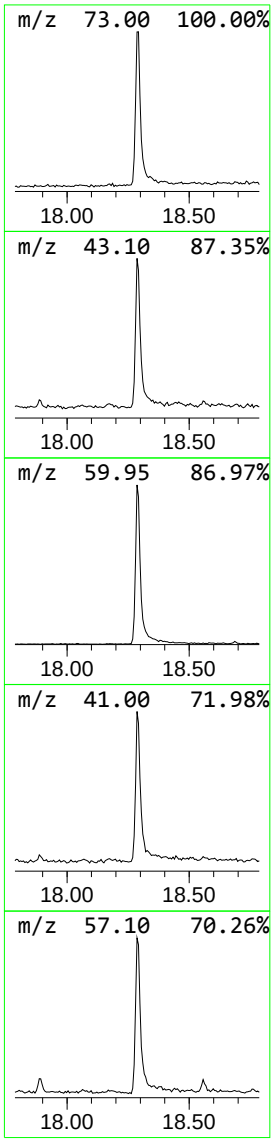
Peak Number 8 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	9.56 ng/ul	1117300	Phenanthrene-d10	17.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91





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Operator : CG/JU
Sample : 02093-14
Misc :
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Instrument :
BNA_P
ClientSampleId :
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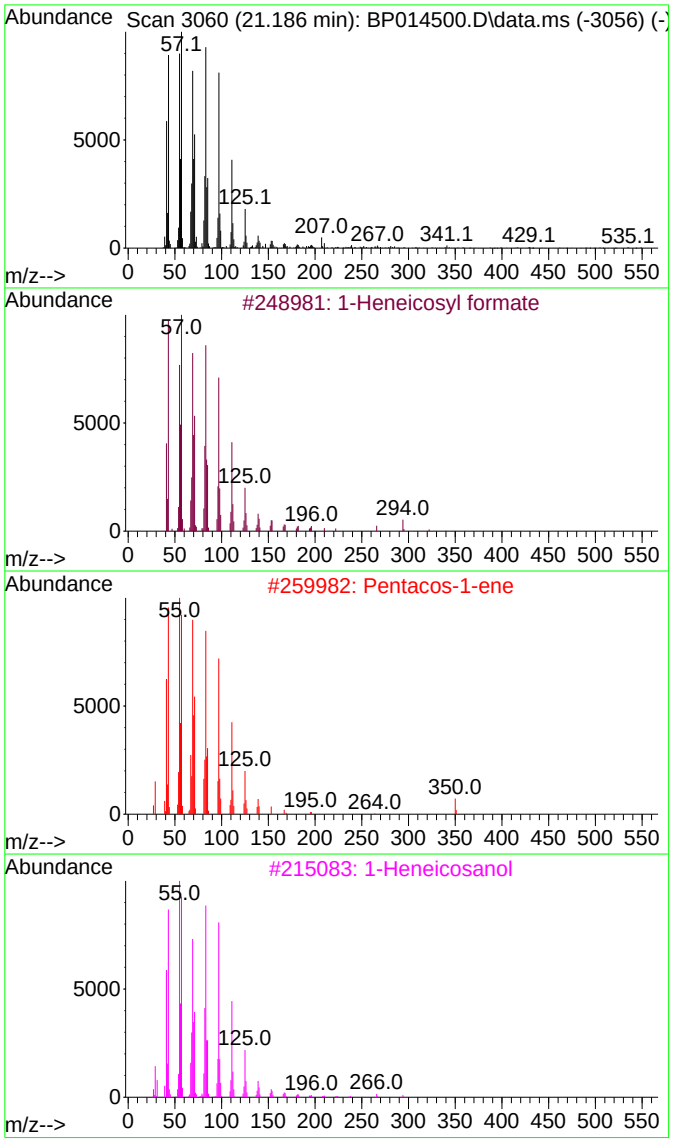
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

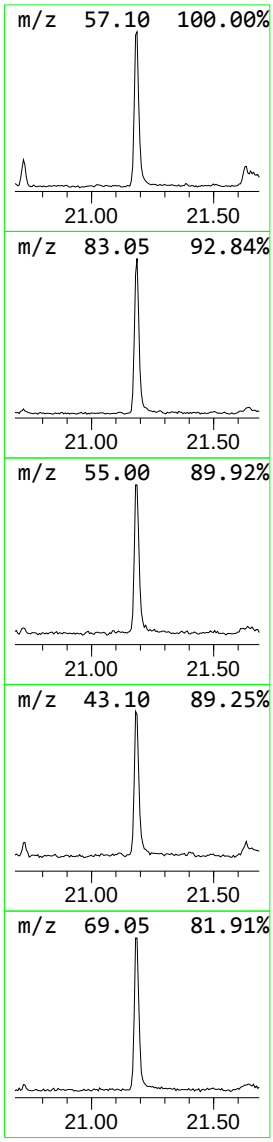
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1-Heneicosyl formate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.186	8.24 ng/ul	1055210	Chrysene-d12	21.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2	Pentacos-1-ene	350	C25H50	016980-85-1	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	Nonacos-1-ene	406	C29H58	018835-35-3	91
5	1-Nonadecene	266	C19H38	018435-45-5	91





Tentatively Identified Compound (LSC) summary

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Data File : BP014500.D
Acq On : 03 Apr 2023 22:36
Operator : CG/JU
Sample : 02093-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB996

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.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
Benzenamine, 3-...	9.010	2.8	ng/ul	115659	1	8.005	813525	20.0
L-Fenchone	9.252	4.6	ng/ul	187865	1	8.005	813525	20.0
Phenol, m-tert-...	12.175	2.0	ng/ul	134033	2	10.828	1321400	20.0
Diethyltoluamide	15.398	6.1	ng/ul	566737	3	14.657	1859800	20.0
Benzophenone	16.022	4.2	ng/ul	390803	3	14.657	1859800	20.0
Benzenesulfonam...	16.181	6.0	ng/ul	705004	4	17.422	2336240	20.0
Benzenesulfonam...	16.692	3.2	ng/ul	376075	4	17.422	2336240	20.0
n-Hexadecanoic ...	18.286	9.6	ng/ul	1117300	4	17.422	2336240	20.0
1-Heneicosyl fo...	21.186	8.2	ng/ul	1055210	5	21.510	2561650	20.0