LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

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_M\Methods\SFAM-EPA-BM091522.M
Title : SVOA CALIBRATION

Signal : TIC: BM036666.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	19	rVB	4449662	5610668	100.00%	7.494%
2	3.393	49	52	54	rBV2	24814	26573	0.47%	0.035%
3	3.428	54	58	69	rVB	590869	759848	13.54%	1.015%
4	3.681	97	101	107	rBV3	54274	75760	1.35%	0.101%
5	3.828	123	126	139	rBV	193125	305922	5.45%	0.409%
6	4.063	161	166	170	rVB	26135	37324	0.67%	0.050%
7	4.157	177	182	189	rBV	26101	40735	0.73%	0.054%
8	4.234	191	195	200	rVV	136113	178169	3.18%	0.238%
9	4.281	200	203	209	rVB2	13459	19540	0.35%	0.026%
10	4.610	252	259	269	rVB	135235	227974	4.06%	0.304%
11	5.104	335	343	355	rBV	72862	123304	2.20%	0.165%
12	5.304	371	377	383	rVB	34540	54089	0.96%	0.072%
13	5.581	419	424	431	rVB9	8964	18166	0.32%	0.024%
14	5.793	456	460	466	rBV9	8964	17556	0.31%	0.023%
15	5.940	481	485	490	rVB6	9763	14543	0.26%	0.019%
16	6.087	506	510	516	rVB7	9499	16381	0.29%	0.022%
17	6.281	539	543	547	rVB	18417	28605	0.51%	0.038%
18	6.598	593	597	604	rBV6	13172	24327	0.43%	0.032%
19	6.722	613	618	623	rVB9	8199	16043	0.29%	0.021%
20	6.998	661	665	669	rVB4	8329	15458	0.28%	0.021%
21	7.175	687	695	707	rVB	259128	438325	7.81%	0.585%
22	7.345	717	724	733	rBV	888642	1347173	24.01%	1.799%
23	7.463	739	744	751	rVB6	8866	17478	0.31%	0.023%
24	7.545	751	758	768	rBV	864360	1331069	23.72%	1.778%
25	8.016	829	838	845	rVV	1168412	1843979	32.87%	2.463%
26	8.098	847	852	860	rVB4	26233	61294	1.09%	0.082%
27	8.486	914	918	925	rVB8	10157	17576	0.31%	0.023%
28	8.557	925	930	932	rBV5	8163	13741	0.24%	0.018%
29	8.728	946	959	966	rBV	770262	1235035	22.01%	1.650%
30	9.010	1000	1007	1015	rBV	863988	1407215	25.08%	1.880%
31	9.128	1022	1027	1031	rBV2	37594	54945	0.98%	0.073%
32	9.186	1031	1037	1044	rVB	988645	1617099	28.82%	2.160%
33	9.292	1053	1055	1062	rVB4	15285	25627	0.46%	0.034%
34	9.398	1068	1073	1079	rVB9	21759	45504	0.81%	0.061%
35	9.545	1092	1098	1104	rBV5	36804	75538	1.35%	0.101%
36	9.622	1106	1111	1115	rVV3	36479	77674	1.38%	0.104%

LSC Area Percent Report

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Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

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_M\Methods\SFAM-EPA-BM091522.M
Title : SVOA CALIBRATION

37	9.663	1115	1118	1124	rVB3	32154	58427	1.04%	0.078%
38	9.775	1133	1137	1142	rBV6	12542	16607	0.30%	0.022%
39	9.833	1143	1147	1153	rVB8	11129	20055	0.36%	0.027%
40	9.910	1153	1160	1170	rBV	736779	1198324	21.36%	1.601%
41	10.057	1180	1185	1190	rVB7	21028	39687	0.71%	0.053%
42	10.139	1195	1199	1204	rVB6	13134	20039	0.36%	0.027%
43	10.233	1212	1215	1221	rVB8	8029	13815	0.25%	0.018%
44	10.322	1225	1230	1233	rBV4	11965	20224	0.36%	0.027%
45	10.386	1236	1241	1245	rVV	38223	72165	1.29%	0.096%
46	10.445	1245	1251	1272	rVB	1314013	2299621	40.99%	3.071%
47	10.610	1277	1279	1290	rVV	14274	34304	0.61%	0.046%
48	10.757	1298	1304	1309	rVV	39680	80752	1.44%	0.108%
49	10.833	1309	1317	1324	rVV	1646310	2779520	49.54%	3.712%
50	10.886	1325	1326	1331	rVV5	13771	14973	0.27%	0.020%
51	10.969	1331	1340	1354	rVV	716133	1292322	23.03%	1.726%
52	11.080	1356	1359	1364	rVV7	16659	30547	0.54%	0.041%
53	11.133	1364	1368	1374	rVV8	14575	34968	0.62%	0.047%
54	11.192	1374	1378	1382	rVB7	17972	29563	0.53%	0.039%
55	11.422	1411	1417	1421	rVB9	19122	35206	0.63%	0.047%
56	11.504	1425	1431	1433	rBV7	8456	16662	0.30%	0.022%
57	11.527	1433	1435	1441	rVB5	15318	22265	0.40%	0.030%
58	11.616	1443	1450	1454	rBV6	27110	59653	1.06%	0.080%
59	11.798	1477	1481	1487	rBV6	14965	22877	0.41%	0.031%
60	12.033	1517	1521	1526	rVB2	35534	53432	0.95%	0.071%
61	12.116	1530	1535	1539	rVV6	31597	59668	1.06%	0.080%
62	12.163	1540	1543	1558	rVB8	42565	95470	1.70%	0.128%
63	12.327	1564	1571	1578	rBV	427628	706849	12.60%	0.944%
64	12.427	1582	1588	1603	rVB3	98886	280437	5.00%	0.375%
65	12.545	1603	1608	1613	rBV5	23185	43208	0.77%	0.058%
66	12.639	1618	1624	1630	rVB	159545	258501	4.61%	0.345%
67	12.780	1644	1648	1650	rBV4	17415	23021	0.41%	0.031%
68	12.816	1651	1654	1660	rVB3	28878	48676	0.87%	0.065%
69	13.033	1682	1691	1695	rBV3	30160	74405	1.33%	0.099%
70	13.204	1716	1720	1723	rBV6	19168	31406	0.56%	0.042%
71	13.380	1747	1750	1754	rBV6	16203	22719	0.40%	0.030%
72	13.551	1776	1779	1785	rVB4	28477	42434	0.76%	0.057%
73	13.692	1800	1803	1808	rBV	22136	33168	0.59%	0.044%
74	13.745	1808	1812	1816	rVB2	28218	37399	0.67%	0.050%
75	14.057	1859	1865	1875	rVB	2616139	3483858	62.09%	4.653%
76	14.198	1886	1889	1895	rVB8	24041	33203	0.59%	0.044%

LSC Area Percent Report

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Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

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_M\Methods\SFAM-EPA-BM091522.M
Title : SVOA CALIBRATION

77	14.339	1907	1913	1926	rBV	2979318	4218719	75.19%	5.635%
78	14.445	1927	1931	1939	rVB	67633	113848	2.03%	0.152%
79	14.568	1947	1952	1957	rVB2	171769	239160	4.26%	0.319%
80	14.645	1959	1965	1976	rVB	2411384	3527874	62.88%	4.712%
81	14.833	1992	1997	2006	rBV	197242	299469	5.34%	0.400%
82	15.386	2086	2091	2096	rBV	151999	237905	4.24%	0.318%
83	15.633	2127	2133	2142	rBV	3811736	5296975	94.41%	7.075%
84	15.739	2146	2151	2160	rVV	813273	1168982	20.83%	1.561%
85	15.862	2170	2172	2183	rVB7	47466	103811	1.85%	0.139%
86	16.145	2214	2220	2229	rBV	839679	1202091	21.43%	1.606%
87	16.468	2272	2275	2280	rBV2	59437	75302	1.34%	0.101%
88	16.651	2302	2306	2311	rVB	477992	617576	11.01%	0.825%
89	17.380	2424	2430	2438	rBV	2854838	3968480	70.73%	5.300%
90	17.480	2441	2447	2456	rVV	3920658	5433796	96.85%	7.257%
91	18.274	2578	2582	2592	rBV	405033	649897	11.58%	0.868%
92	19.051	2711	2714	2719	rVB2	109618	142819	2.55%	0.191%
93	19.545	2795	2798	2803	rBV	222205	253694	4.52%	0.339%
94	19.762	2830	2835	2840	rBV	4287099	5605545	99.91%	7.487%
95	19.809	2840	2843	2850	rVB	297072	459404	8.19%	0.614%
96	20.756	3001	3004	3008	rBV	313207	348425	6.21%	0.465%
97	21.256	3086	3089	3094	rVB	463907	496121	8.84%	0.663%
98	21.544	3134	3138	3147	rVB	2487233	3220698	57.40%	4.302%
99	23.833	3521	3527	3534	rVB	1988910	3702736	65.99%	4.945%
100	23.991	3548	3554	3562	rVB2	1335573	2727693	48.62%	3.643%

Sum of corrected areas: 74871707

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
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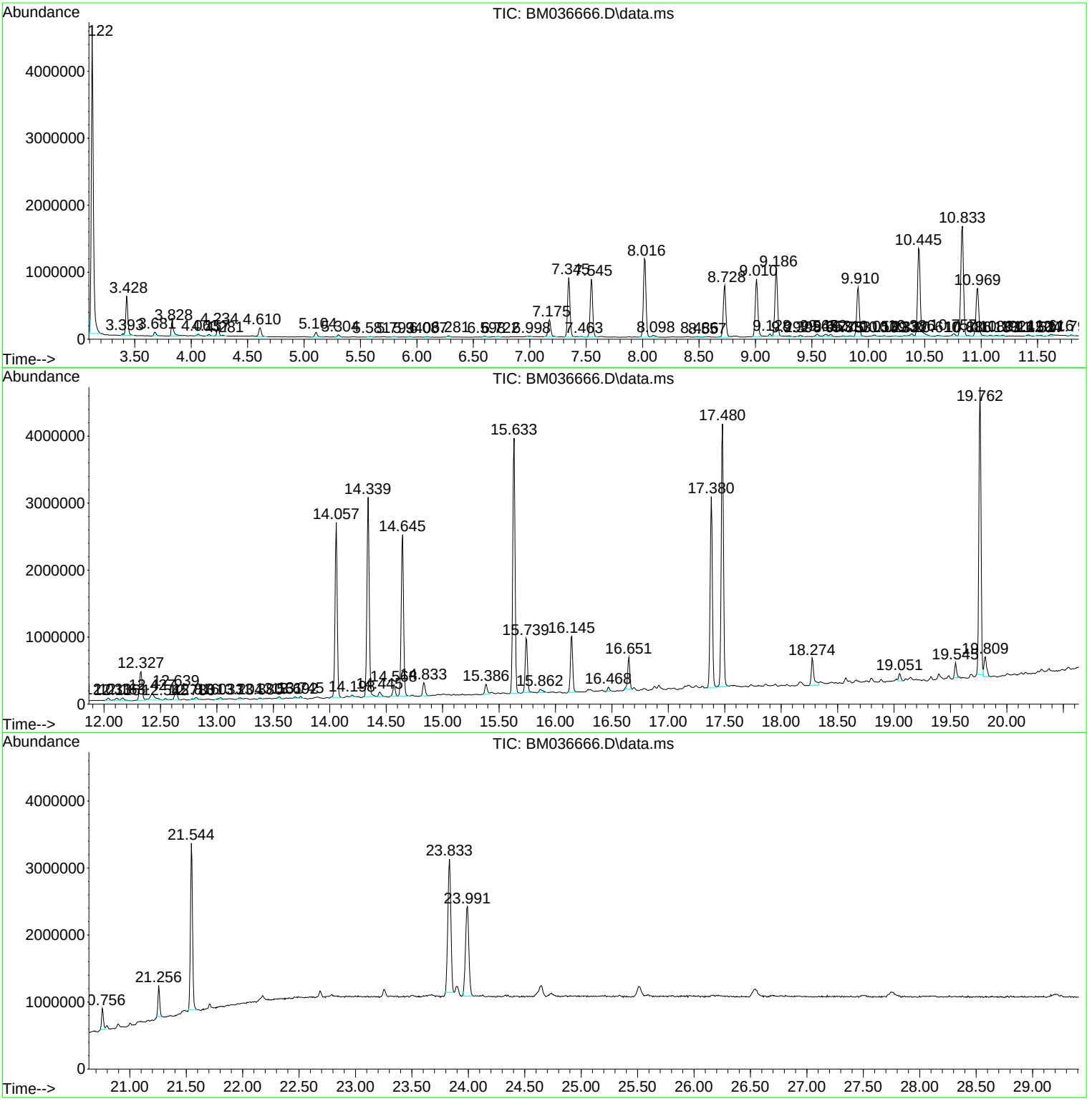
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
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Sample : N4595-06DL 2X
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Instrument :
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ClientSampleId :
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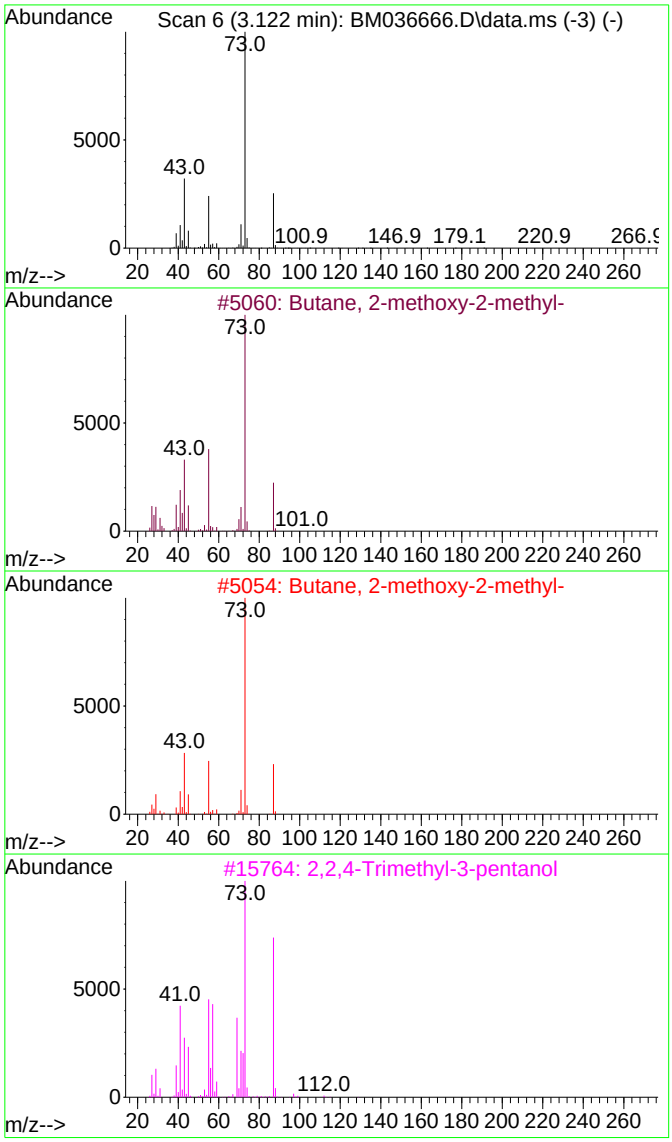
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

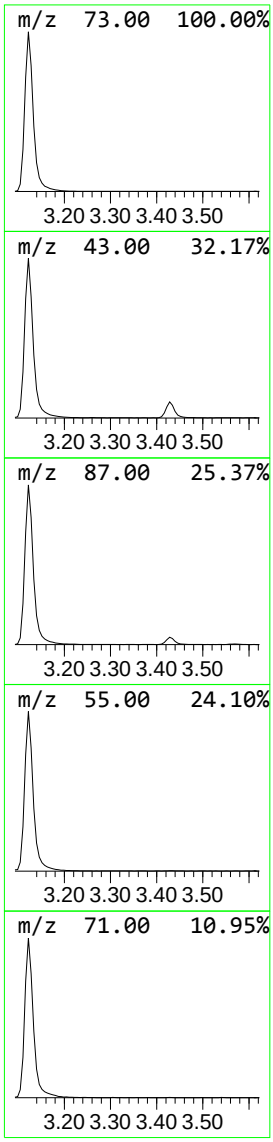
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	60.85 ng/ul	5610670	1,4-Dichlorobenzene-d4	8.022

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

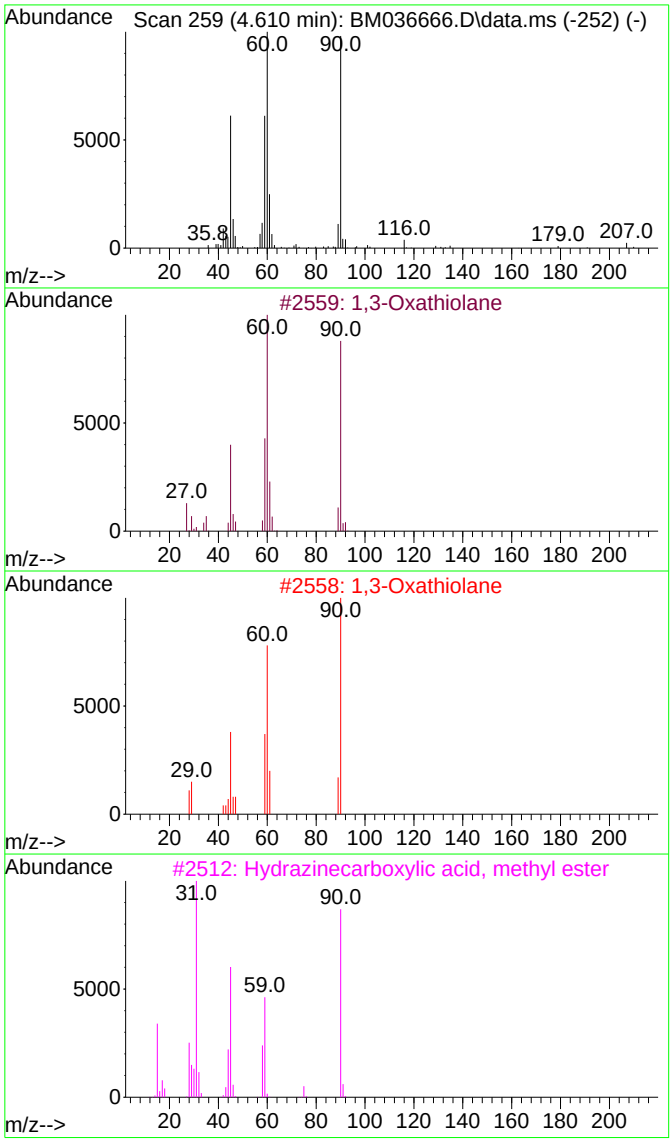
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

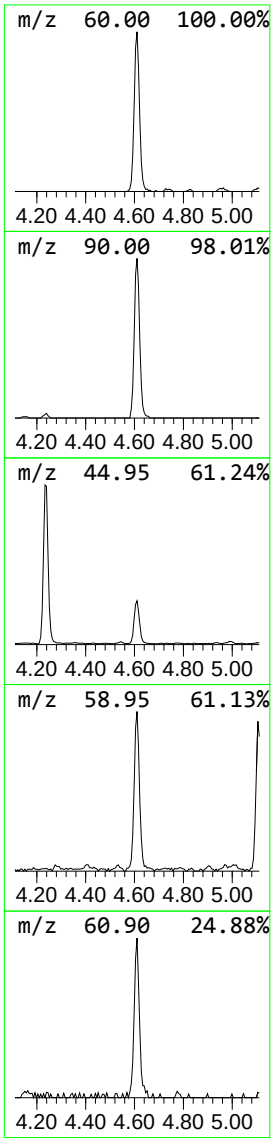
****

Peak Number 2 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.610	2.47 ng/ul	227974	1,4-Dichlorobenzene-d4	8.022

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	94
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
3	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	45
4	3-Thietanol	90	C3H6OS	010304-16-2	40
5	3-Thietanol	90	C3H6OS	010304-16-2	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

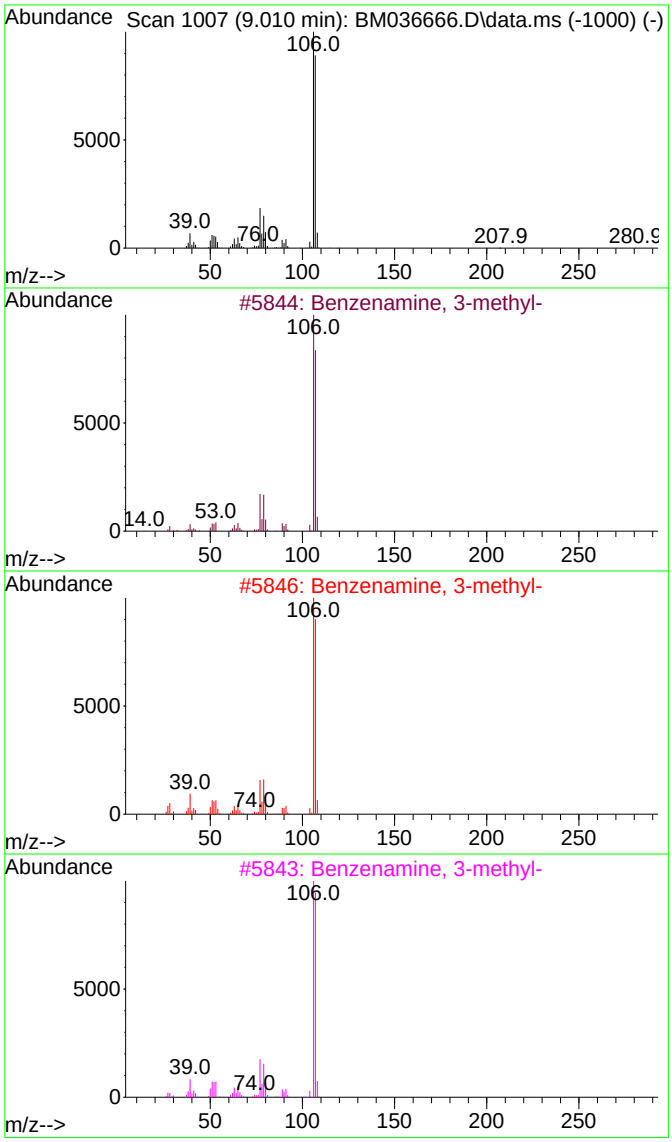
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

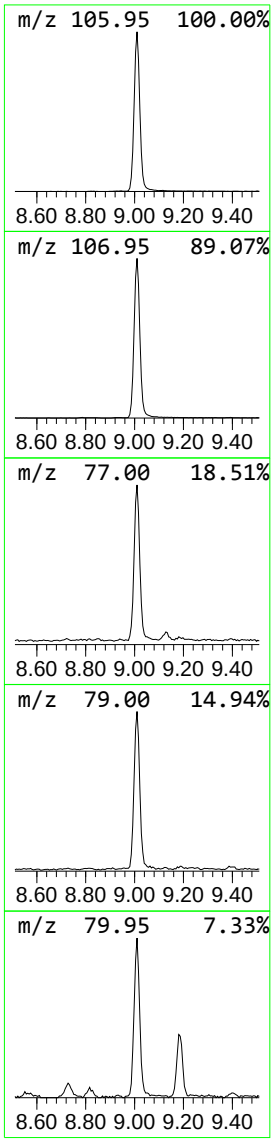
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzenamine, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.010	15.26 ng/ul	1407220	1,4-Dichlorobenzene-d4	8.022

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	o-Toluidine	107	C7H9N	000095-53-4	94





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Instrument :
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ClientSampleId :
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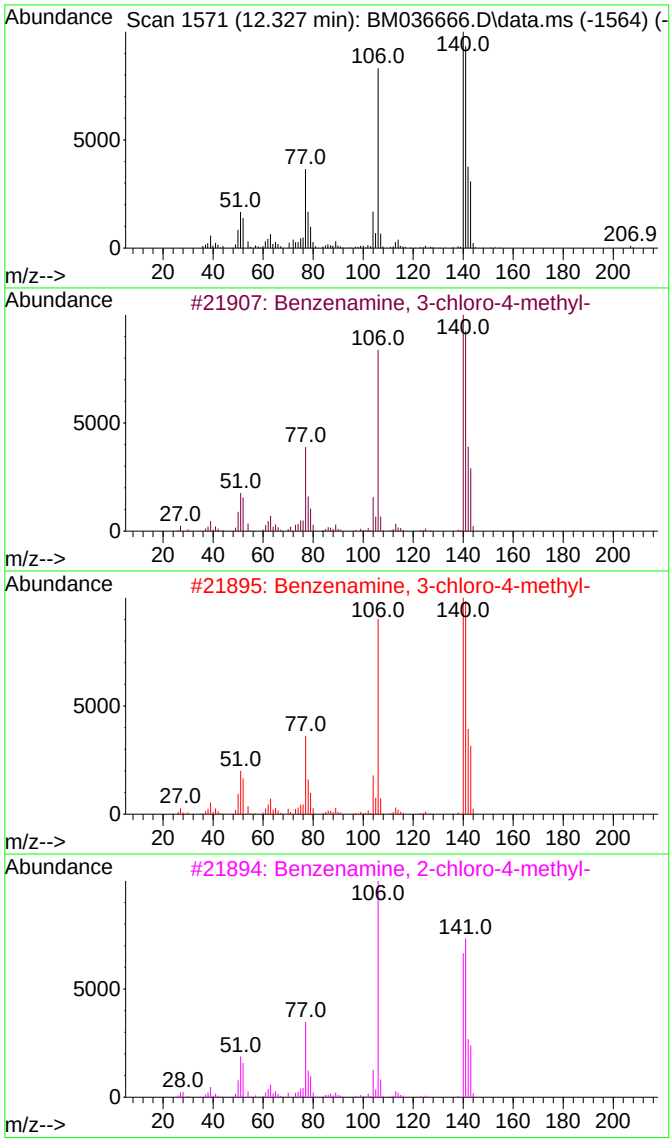
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

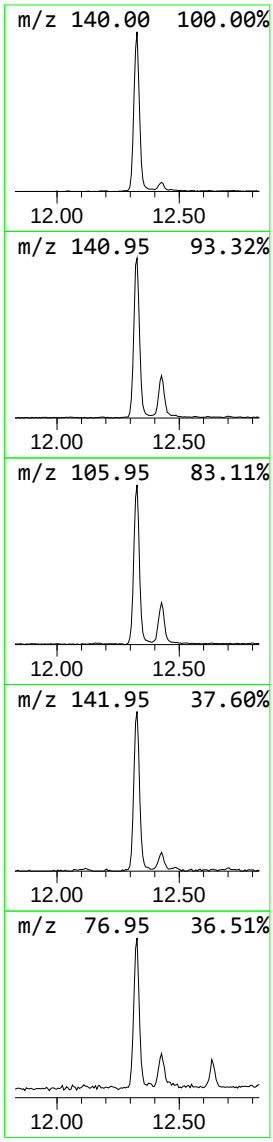
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzenamine, 3-chloro-4-met... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.327	5.09 ng/ul	706849	Naphthalene-d8	10.833

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
3	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	95
4	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	95
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

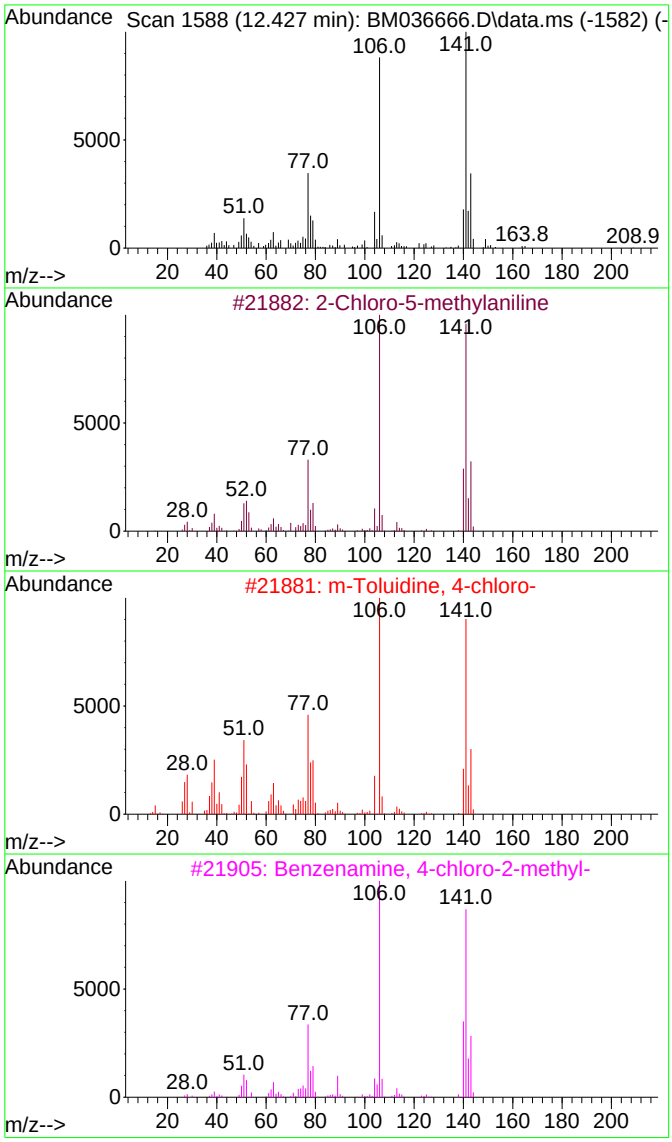
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

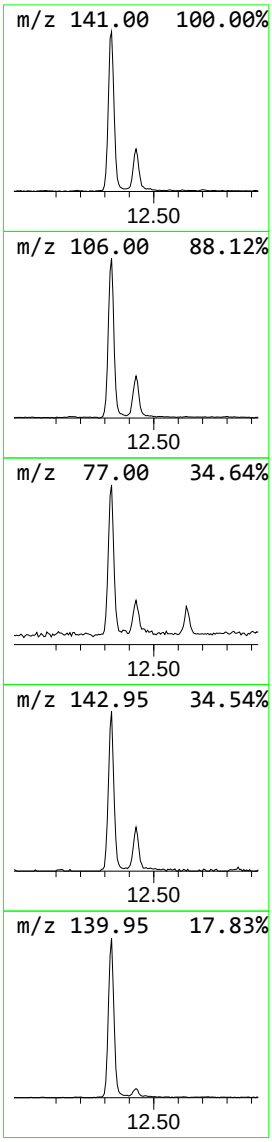
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Chloro-5-methylaniline Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.427	2.02 ng/ul	280437	Naphthalene-d8	10.833

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-methylaniline	141	C7H8ClN	000095-81-8	93
2	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	91
3	Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	90
4	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	87
5	Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	87





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Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

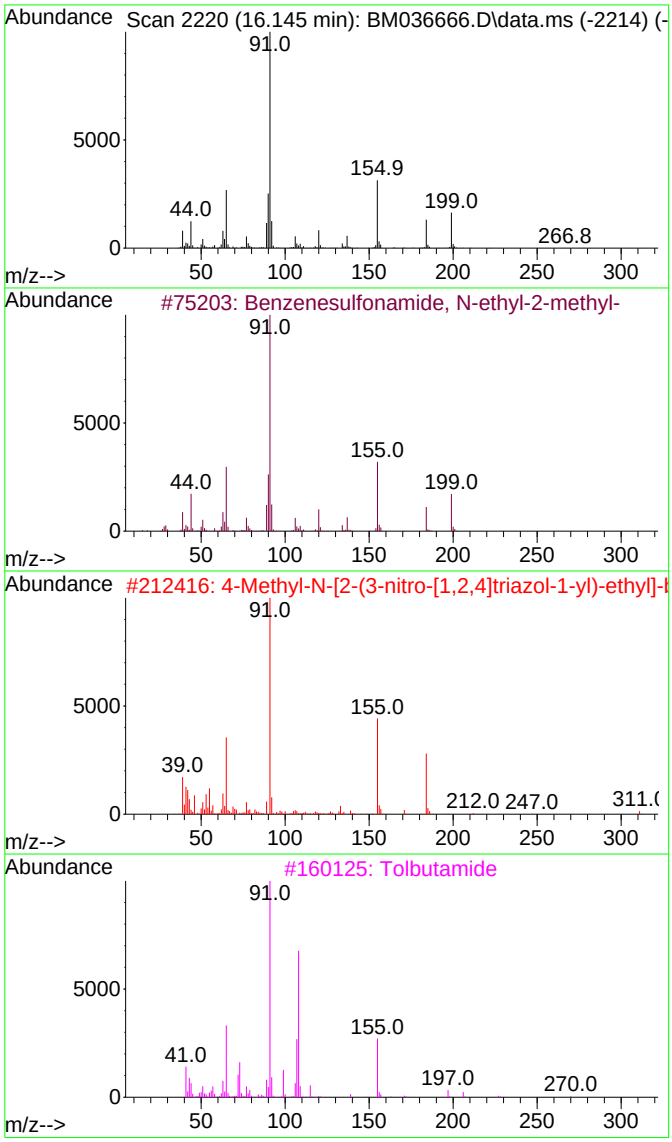
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

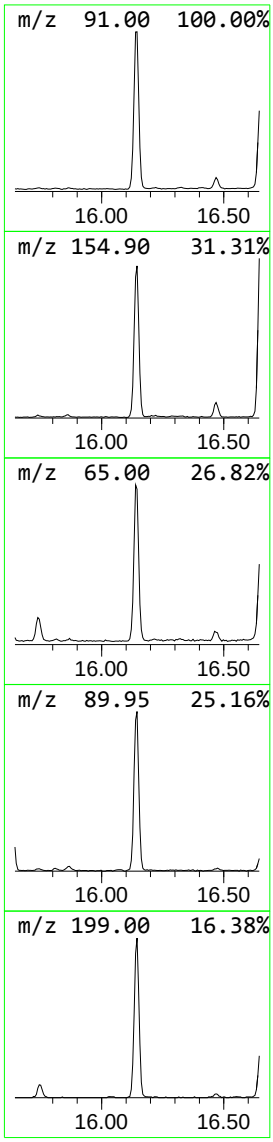
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	6.06 ng/ul	1202090	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	50
3	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
4	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50
5	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	49





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Data File : BM036666.D
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Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

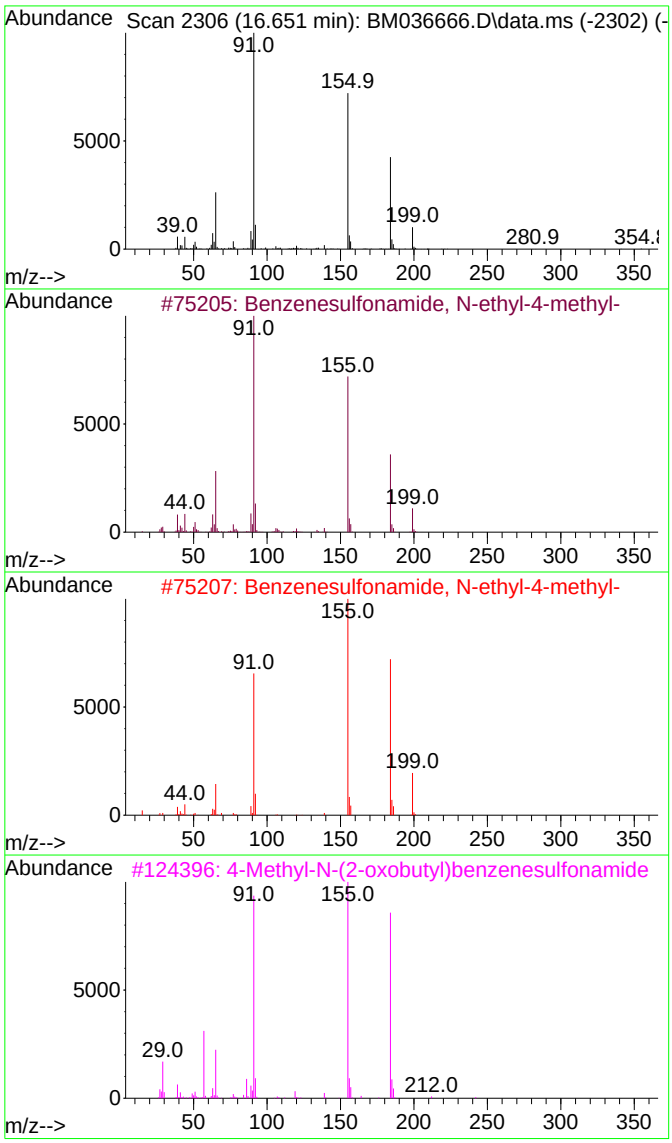
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

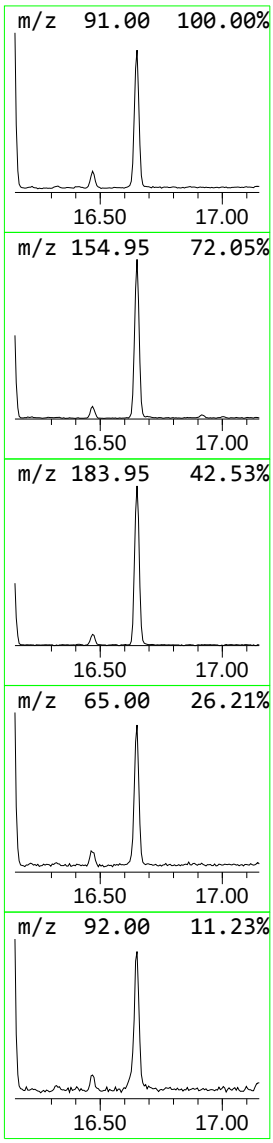
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	3.11 ng/ul	617576	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	90
3	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
4	Benzene, 4-(ethylsulfonyl)-1-met...	184	C9H12O2S	007569-34-8	80
5	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78





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Data File : BM036666.D
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Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
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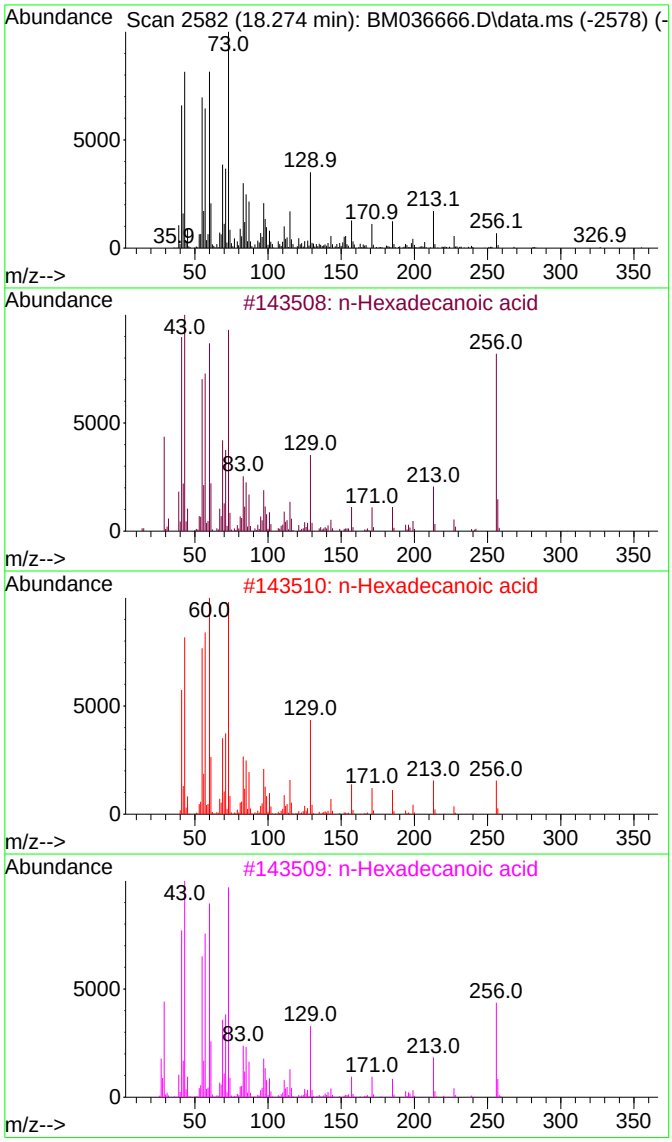
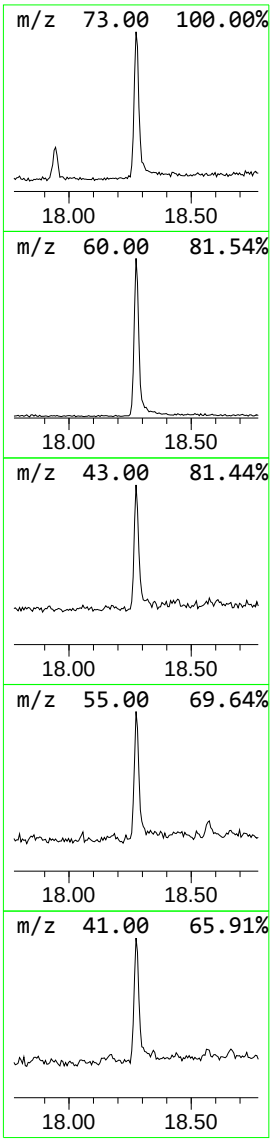
****

Peak Number 8 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.274	3.28 ng/ul	649897	Phenanthrene-d10	17.380

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	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	95





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Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

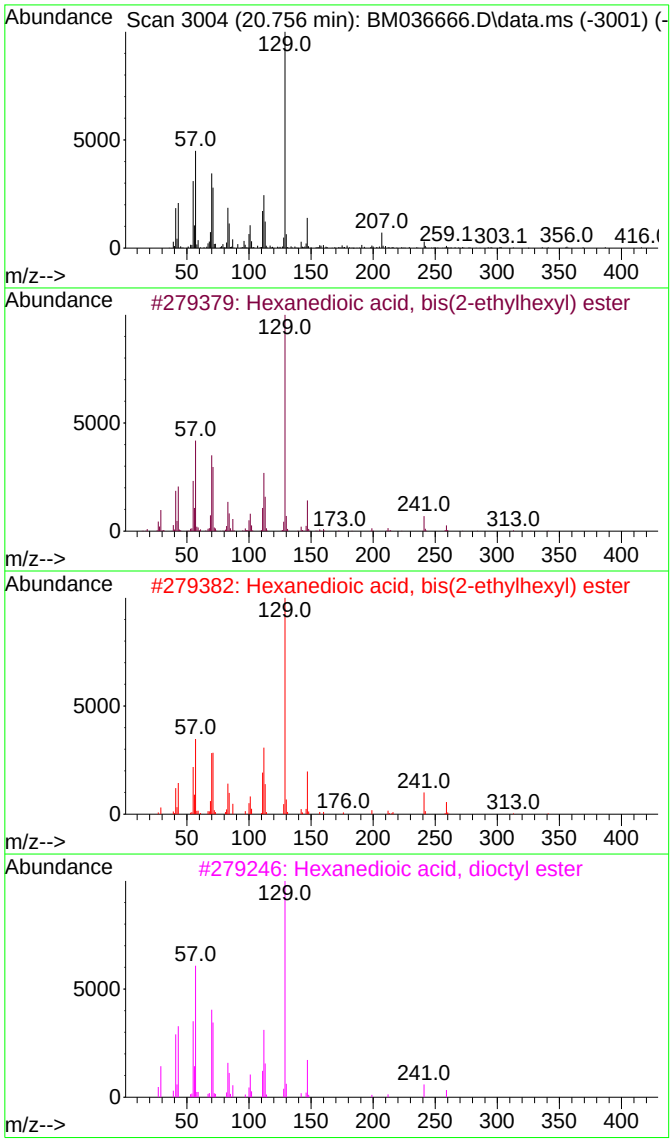
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

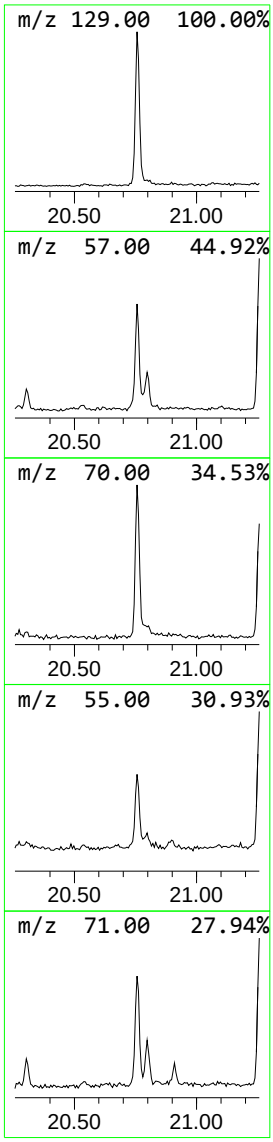
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Hexanedioic acid, bis(2-eth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.756	2.16 ng/ul	348425	Chrysene-d12	21.544

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
3	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	91
4	Diisooctyl adipate	370	C22H42O4	001330-86-5	78
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76





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Data File : BM036666.D
Acq On : 22 Sep 2022 17:14
Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H4DL

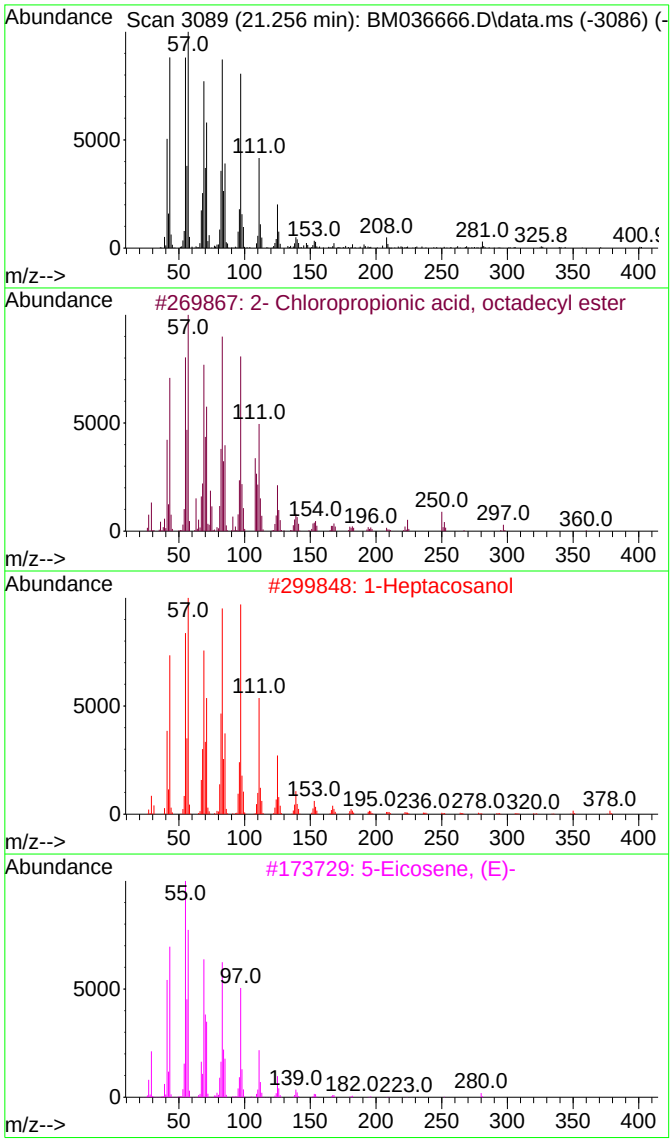
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

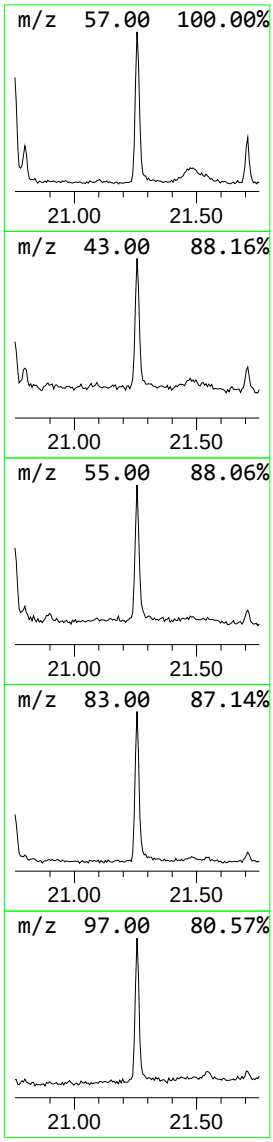
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2- Chloropropionic acid, oc... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.256	3.08 ng/ul	496121	Chrysene-d12	21.544

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2	1-Heptacosanol	396	C27H56O	002004-39-9	94
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	1-Hexacosene	364	C26H52	018835-33-1	93





Tentatively Identified Compound (LSC) summary

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Operator : CG/JU
Sample : N4595-06DL 2X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.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...	3.122	60.9	ng/ul	5610670	1	8.022	1843980	20.0
1,3-Oxathiolane	4.610	2.5	ng/ul	227974	1	8.022	1843980	20.0
Benzenamine, 3-...	9.010	15.3	ng/ul	1407220	1	8.022	1843980	20.0
Benzenamine, 3-...	12.327	5.1	ng/ul	706849	2	10.833	2779520	20.0
2-Chloro-5-meth...	12.427	2.0	ng/ul	280437	2	10.833	2779520	20.0
Benzenesulfonam...	16.145	6.1	ng/ul	1202090	4	17.380	3968480	20.0
Benzenesulfonam...	16.651	3.1	ng/ul	617576	4	17.380	3968480	20.0
n-Hexadecanoic ...	18.274	3.3	ng/ul	649897	4	17.380	3968480	20.0
Hexanedioic aci...	20.756	2.2	ng/ul	348425	5	21.544	3220700	20.0
2- Chloropropio...	21.256	3.1	ng/ul	496121	5	21.544	3220700	20.0