LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G9DL

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: BM036664.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	18	rVB	1889699	2283668	80.28%	6.093%
2	3.393	49	52	53	rBV2	11054	10654	0.37%	0.028%
3	3.428	53	58	72	rVB	734709	933701	32.82%	2.491%
4	3.675	96	100	107	rBV	69823	96089	3.38%	0.256%
5	3.834	126	127	138	rBV	66461	121155	4.26%	0.323%
6	4.063	162	166	172	rVB3	13244	25516	0.90%	0.068%
7	4.152	176	181	187	rBV	19557	31410	1.10%	0.084%
8	4.234	191	195	205	rVB	122046	169628	5.96%	0.453%
9	4.540	242	247	251	rVV7	10597	19903	0.70%	0.053%
10	4.610	254	259	266	rVB	129171	196730	6.92%	0.525%
11	5.593	418	426	430	rBV8	10956	26663	0.94%	0.071%
12	5.940	480	485	488	rVV4	11819	17179	0.60%	0.046%
13	6.093	505	511	516	rBV10	6310	9990	0.35%	0.027%
14	6.281	534	543	549	rBV	36679	66184	2.33%	0.177%
15	6.522	577	584	588	rBV	21534	34770	1.22%	0.093%
16	6.598	592	597	603	rVB8	10234	20929	0.74%	0.056%
17	6.716	606	617	625	rVV2	88394	151694	5.33%	0.405%
18	6.993	656	664	670	rBV2	10326	26401	0.93%	0.070%
19	7.175	686	695	701	rBV	93365	169342	5.95%	0.452%
20	7.345	717	724	734	rBV	361548	583930	20.53%	1.558%
21	7.445	738	741	748	rVV4	10911	19223	0.68%	0.051%
22	7.545	752	758	766	rVV	335282	530372	18.64%	1.415%
23	7.645	769	775	778	rVB7	10712	19909	0.70%	0.053%
24	7.875	806	814	820	rBV10	7581	17044	0.60%	0.045%
25	8.016	832	838	845	rVV	897388	1449729	50.96%	3.868%
26	8.092	845	851	860	rVB4	36692	81208	2.85%	0.217%
27	8.492	917	919	925	rVB7	8076	10358	0.36%	0.028%
28	8.563	925	931	933	rBV6	9264	18156	0.64%	0.048%
29	8.728	952	959	968	rVV	260902	446883	15.71%	1.192%
30	8.810	968	973	974	rVV5	17488	25059	0.88%	0.067%
31	8.828	974	976	983	rVB4	18331	31205	1.10%	0.083%
32	8.945	989	996	1002	rBV4	8061	21119	0.74%	0.056%
33	9.128	1023	1027	1030	rVV2	20137	30307	1.07%	0.081%
34	9.181	1030	1036	1043	rVV	389172	649693	22.84%	1.733%
35	9.298	1051	1056	1062	rVB7	16393	34776	1.22%	0.093%
36	9.398	1069	1073	1079	rVB9	9614	14174	0.50%	0.038%

LSC Area Percent Report

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Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

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	9.545	1093	1098	1103	rVV3	46473	83239	2.93%	0.222%
38	9.604	1107	1108	1112	rVV4	10262	15342	0.54%	0.041%
39	9.663	1112	1118	1126	rVB4	31765	72908	2.56%	0.195%
40	9.739	1126	1131	1138	rBV9	6073	16927	0.60%	0.045%
41	9.834	1141	1147	1154	rBV10	16880	38280	1.35%	0.102%
42	9.910	1154	1160	1166	rBV	273021	436073	15.33%	1.163%
43	9.957	1166	1168	1172	rVB5	7928	10324	0.36%	0.028%
44	10.022	1174	1179	1183	rBV7	11155	17659	0.62%	0.047%
45	10.139	1194	1199	1202	rVB7	6068	10826	0.38%	0.029%
46	10.186	1202	1207	1208	rBV5	8534	11663	0.41%	0.031%
47	10.322	1225	1230	1233	rBV4	12963	21700	0.76%	0.058%
48	10.381	1233	1240	1245	rVV	86305	161147	5.66%	0.430%
49	10.445	1245	1251	1261	rVV	446442	781282	27.46%	2.084%
50	10.522	1261	1264	1271	rVB4	30405	53261	1.87%	0.142%
51	10.610	1274	1279	1280	rBV5	11166	16272	0.57%	0.043%
52	10.757	1300	1304	1309	rBV2	20237	32775	1.15%	0.087%
53	10.833	1309	1317	1324	rBV	1275802	2136643	75.11%	5.700%
54	10.963	1332	1339	1353	rBV	232276	472638	16.61%	1.261%
55	11.081	1355	1359	1365	rBV9	16827	34905	1.23%	0.093%
56	11.186	1375	1377	1384	rVB6	12316	22703	0.80%	0.061%
57	11.422	1410	1417	1423	rVB9	28577	61412	2.16%	0.164%
58	11.498	1423	1430	1433	rBV7	14044	29329	1.03%	0.078%
59	11.616	1441	1450	1452	rBV10	15397	37037	1.30%	0.099%
60	11.804	1478	1482	1484	rBV4	15044	20521	0.72%	0.055%
61	11.957	1505	1508	1510	rBV4	7023	10134	0.36%	0.027%
62	12.033	1516	1521	1526	rVB	52805	79362	2.79%	0.212%
63	12.092	1526	1531	1532	rBV5	17459	24190	0.85%	0.065%
64	12.163	1540	1543	1554	rVB2	39047	81243	2.86%	0.217%
65	12.328	1564	1571	1577	rBV	184372	305086	10.72%	0.814%
66	12.427	1584	1588	1603	rVB9	52940	178396	6.27%	0.476%
67	12.545	1603	1608	1615	rBV5	28729	62103	2.18%	0.166%
68	12.639	1618	1624	1631	rVV2	120711	211218	7.42%	0.564%
69	12.710	1633	1636	1639	rVB4	15597	20067	0.71%	0.054%
70	12.775	1644	1647	1652	rBV6	16319	23929	0.84%	0.064%
71	13.204	1716	1720	1728	rVB8	34143	64633	2.27%	0.172%
72	13.327	1739	1741	1745	rBV5	9591	11527	0.41%	0.031%
73	13.380	1745	1750	1756	rBV8	32390	73643	2.59%	0.196%
74	13.686	1799	1802	1808	rBV	34717	54915	1.93%	0.147%
75	13.874	1829	1834	1845	rVV	96595	201424	7.08%	0.537%
76	14.057	1859	1865	1876	rVB	903846	1241635	43.65%	3.313%

LSC Area Percent Report

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Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8G9DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Title : SVOA CALIBRATION

77	14.151	1877	1881	1885	rBV	70485	112176	3.94%	0.299%
78	14.339	1907	1913	1923	rBV	953279	1423812	50.05%	3.799%
79	14.439	1927	1930	1943	rVB3	58716	132707	4.66%	0.354%
80	14.569	1947	1952	1957	rVV	217755	299252	10.52%	0.798%
81	14.645	1958	1965	1976	rVB	1841602	2660635	93.53%	7.098%
82	14.833	1994	1997	2003	rVB3	57136	81373	2.86%	0.217%
83	15.380	2086	2090	2095	rBV	146628	229996	8.08%	0.614%
84	15.633	2127	2133	2140	rBV	1229556	1792681	63.02%	4.783%
85	15.739	2146	2151	2156	rVB	235272	329675	11.59%	0.880%
86	16.145	2214	2220	2225	rBV	845094	1203689	42.31%	3.211%
87	16.310	2244	2248	2256	rVB3	216404	336273	11.82%	0.897%
88	16.651	2301	2306	2310	rVB	443761	602560	21.18%	1.608%
89	16.692	2310	2313	2319	rVB3	69620	88439	3.11%	0.236%
90	17.380	2424	2430	2438	rBV	2014608	2844741	100.00%	7.590%
91	17.480	2441	2447	2458	rVB	1213962	1783955	62.71%	4.760%
92	18.274	2579	2582	2590	rBV3	111582	191222	6.72%	0.510%
93	18.568	2630	2632	2636	rBV3	78413	97613	3.43%	0.260%
94	18.798	2668	2671	2676	rVB2	103379	135214	4.75%	0.361%
95	19.051	2711	2714	2718	rVB	113366	132261	4.65%	0.353%
96	19.762	2830	2835	2839	rBV	1283340	1707607	60.03%	4.556%
97	20.756	3001	3004	3010	rBV	332952	358454	12.60%	0.956%
98	21.545	3134	3138	3145	rVB	1721193	2183126	76.74%	5.824%
99	23.833	3522	3527	3534	rVB2	591265	1153893	40.56%	3.079%
100	23.986	3548	3553	3563	rVB2	980163	1967390	69.16%	5.249%

Sum of corrected areas: 37481866

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
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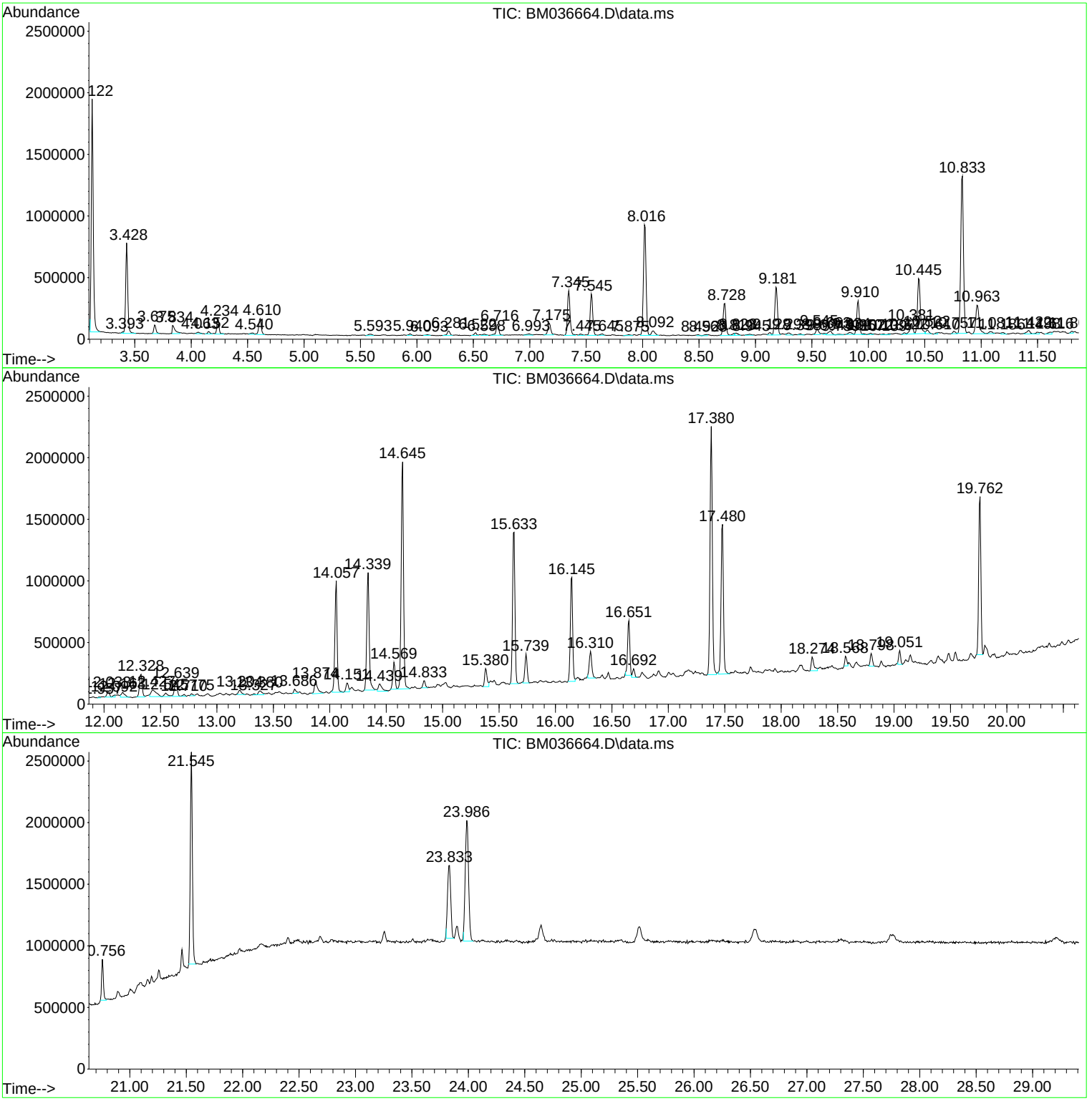
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

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 : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
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Misc :
ALS Vial : 50 Sample Multiplier: 1

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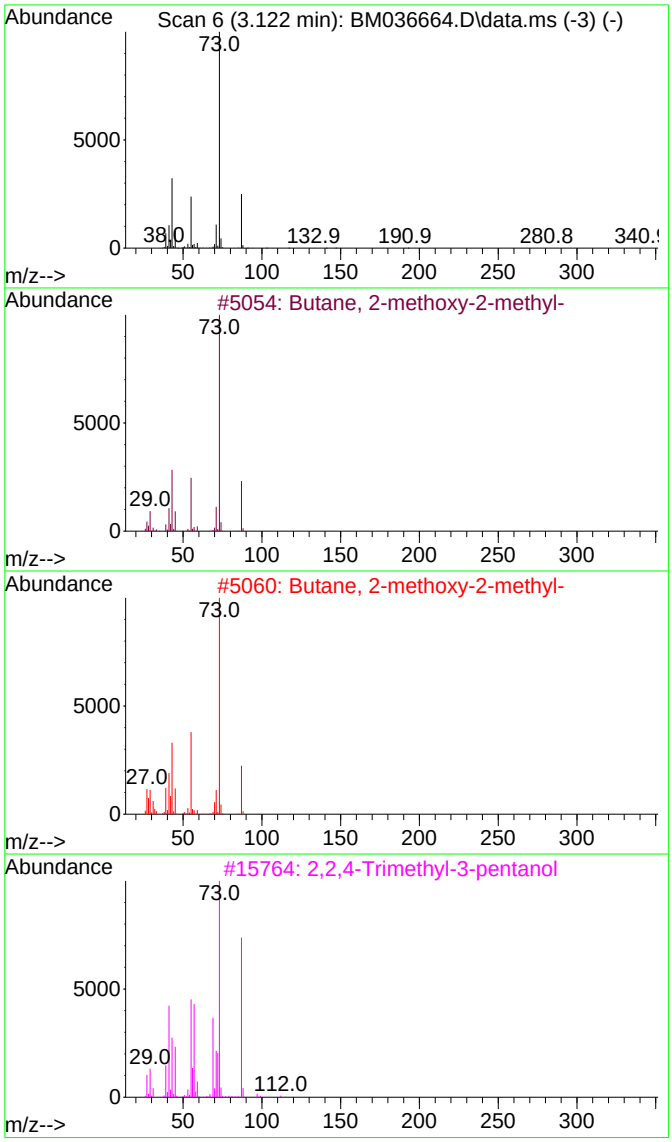
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

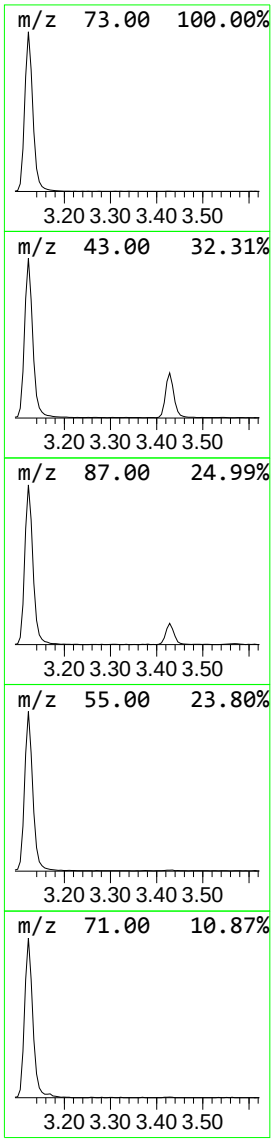
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	31.50 ng/ul	2283670	1,4-Dichlorobenzene-d4	8.016

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	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 : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

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

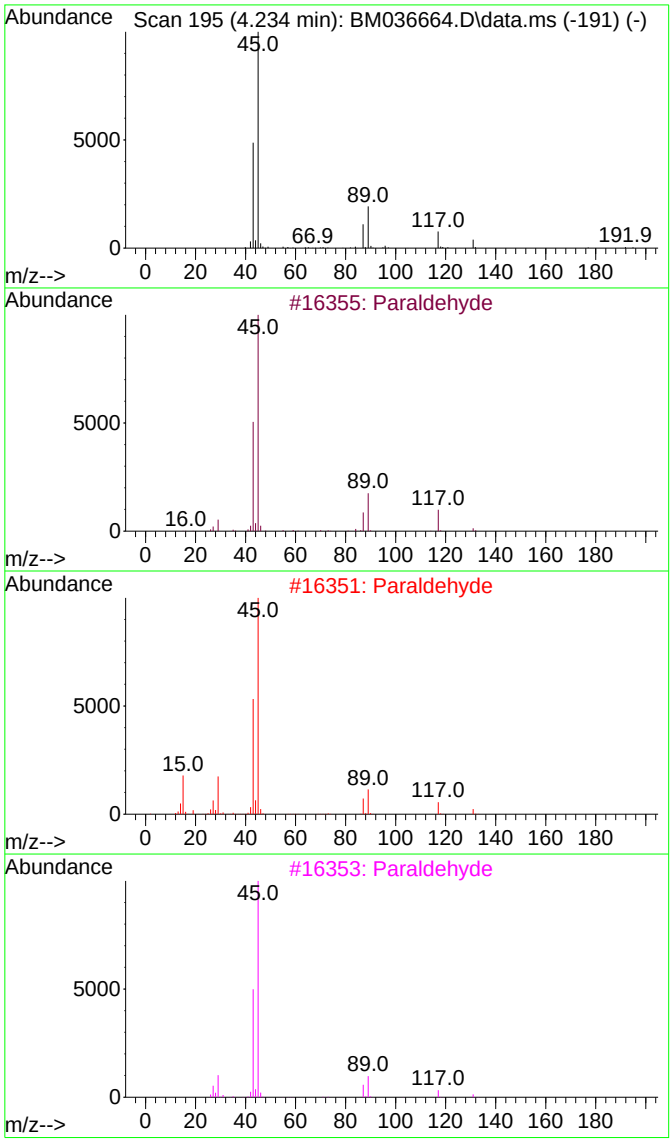
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

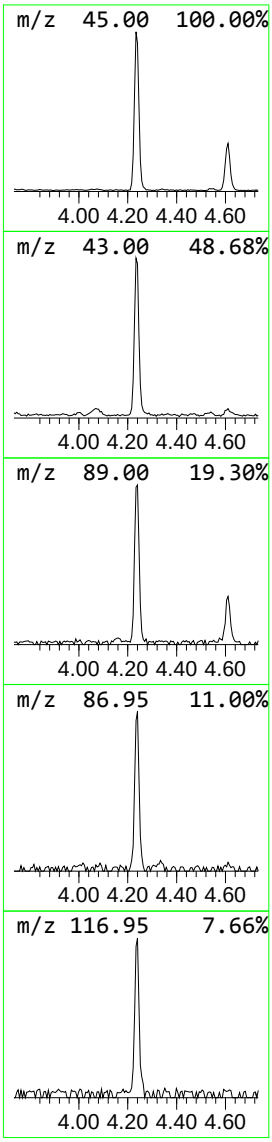
****

Peak Number 2 Paraldehyde Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.234	2.34 ng/ul	169628	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Paraldehyde	132	C6H12O3	000123-63-7	86
2	Paraldehyde	132	C6H12O3	000123-63-7	78
3	Paraldehyde	132	C6H12O3	000123-63-7	64
4	Paraldehyde	132	C6H12O3	000123-63-7	56
5	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
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Acq On : 22 Sep 2022 15:58
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Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 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

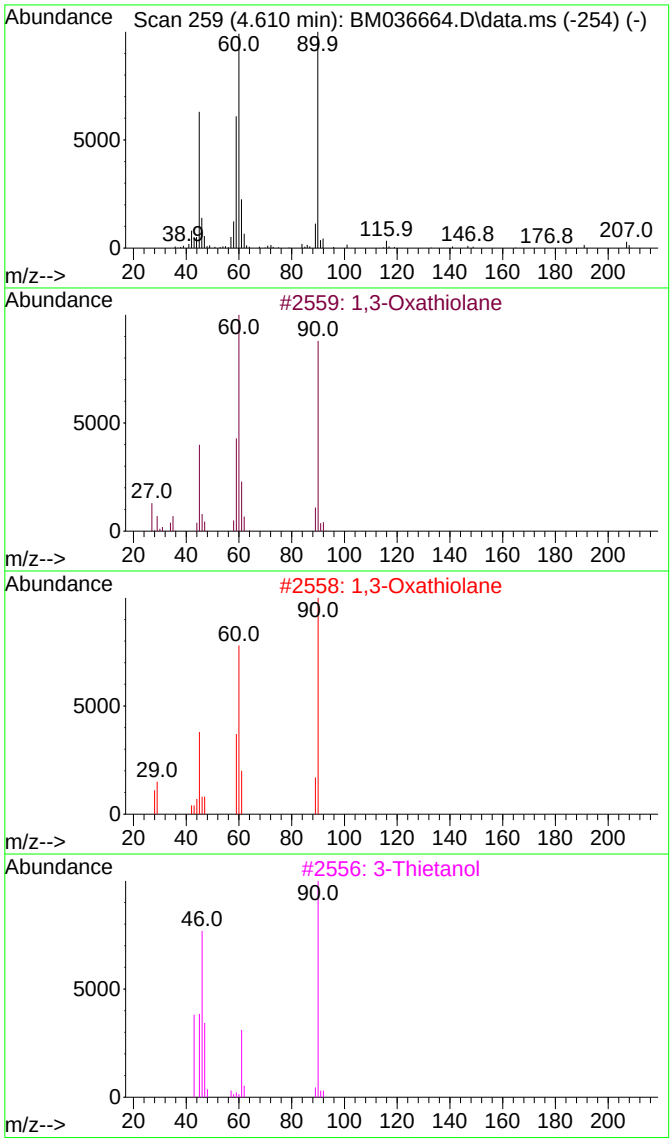
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

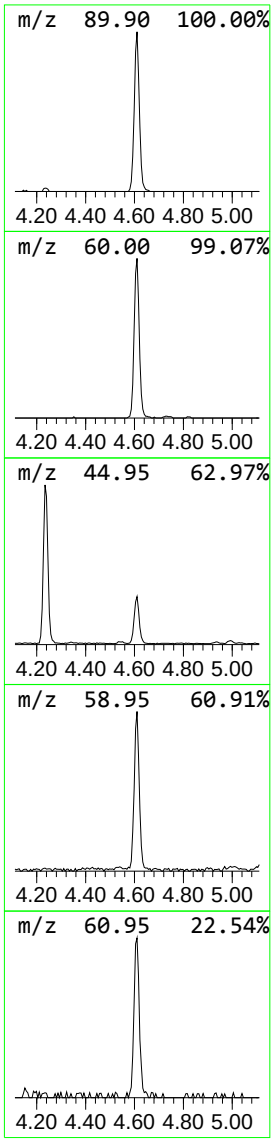
****

Peak Number 3 1,3-Oxathiolane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.610	2.71 ng/ul	196730	1,4-Dichlorobenzene-d4	8.016

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	56
3	3-Thietanol	90	C3H6OS	010304-16-2	53
4	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
5	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50





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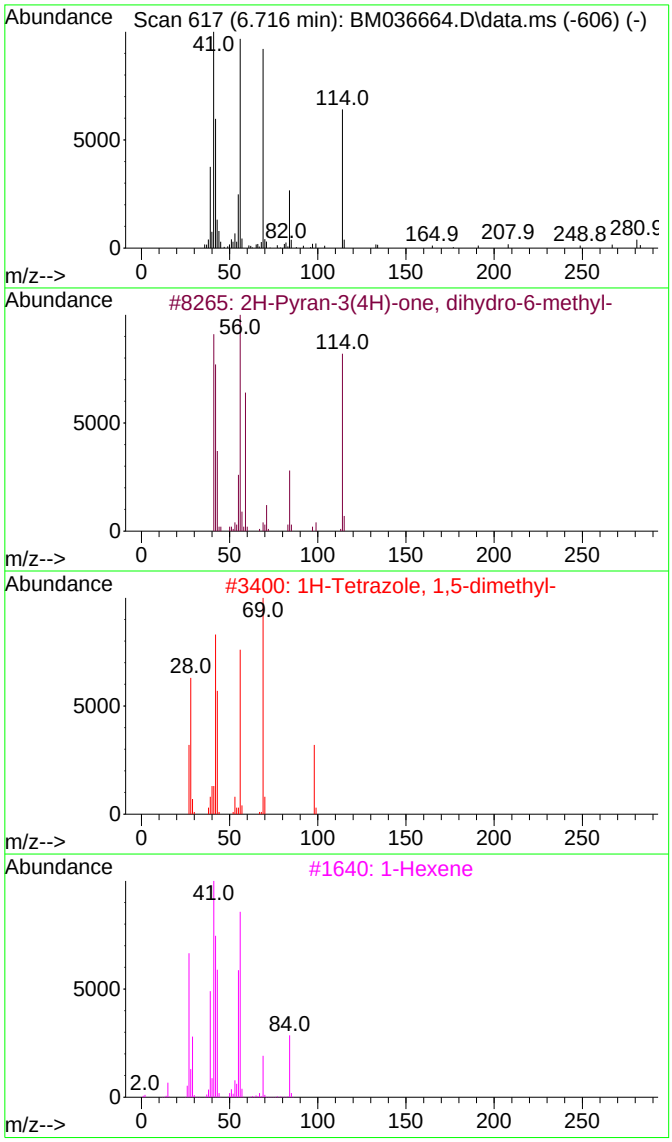
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

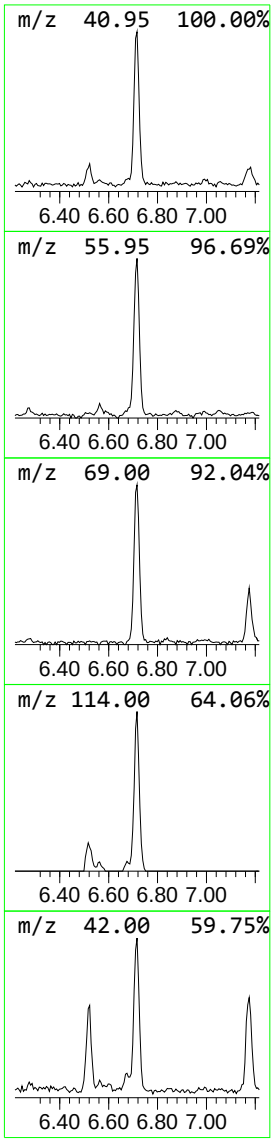
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2H-Pyran-3(4H)-one, dihydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.716	2.09 ng/ul	151694	1,4-Dichlorobenzene-d4	8.016

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran-3(4H)-one, dihydro-6-me...	114	C6H10O2	043152-89-2	62
2	1H-Tetrazole, 1,5-dimethyl-	98	C3H6N4	005144-11-6	50
3	1-Hexene	84	C6H12	000592-41-6	49
4	Pentane, 3-methylene-	84	C6H12	000760-21-4	43
5	2-Pentene, 4-methyl-, (E)-	84	C6H12	000674-76-0	38





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Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

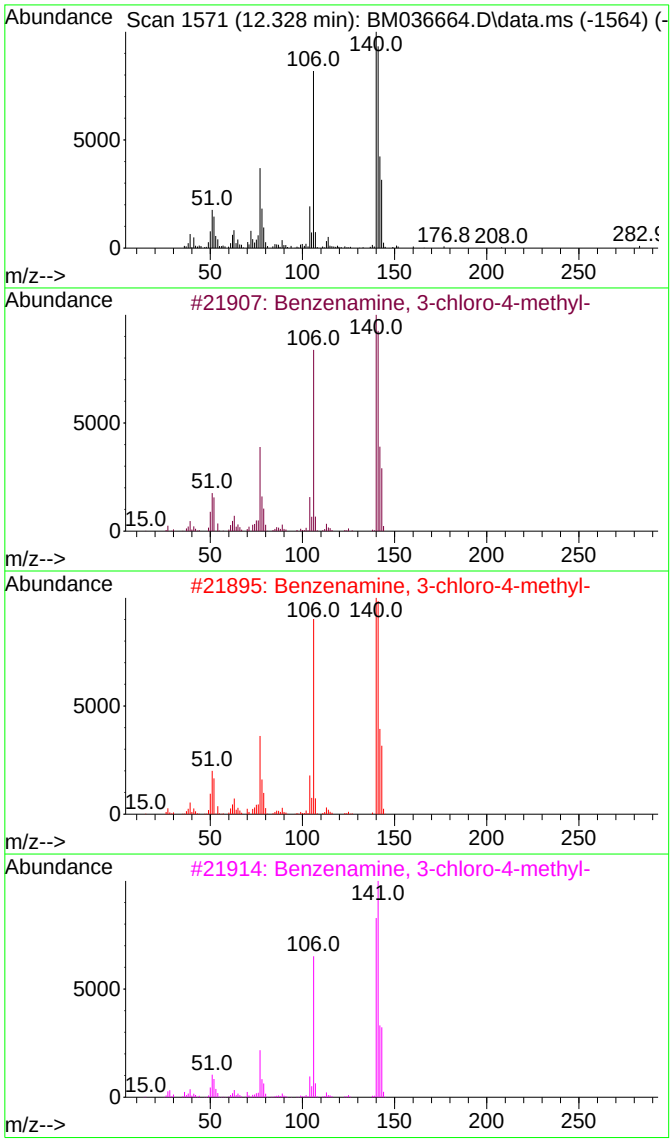
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

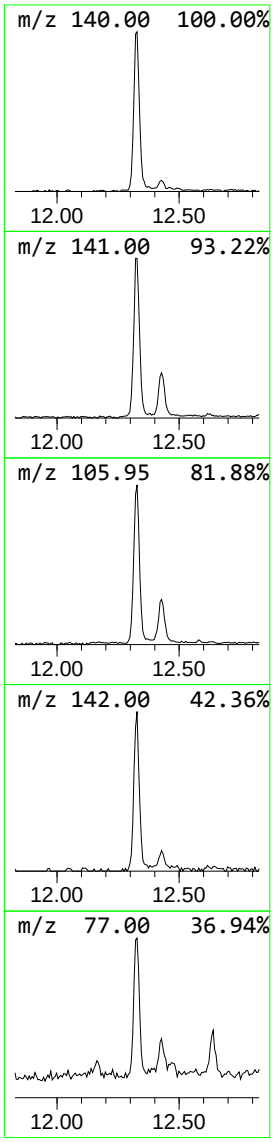
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzenamine, 3-chloro-4-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.328	2.86 ng/ul	305086	Naphthalene-d8	10.834

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, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	95
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	93
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	90





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Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

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

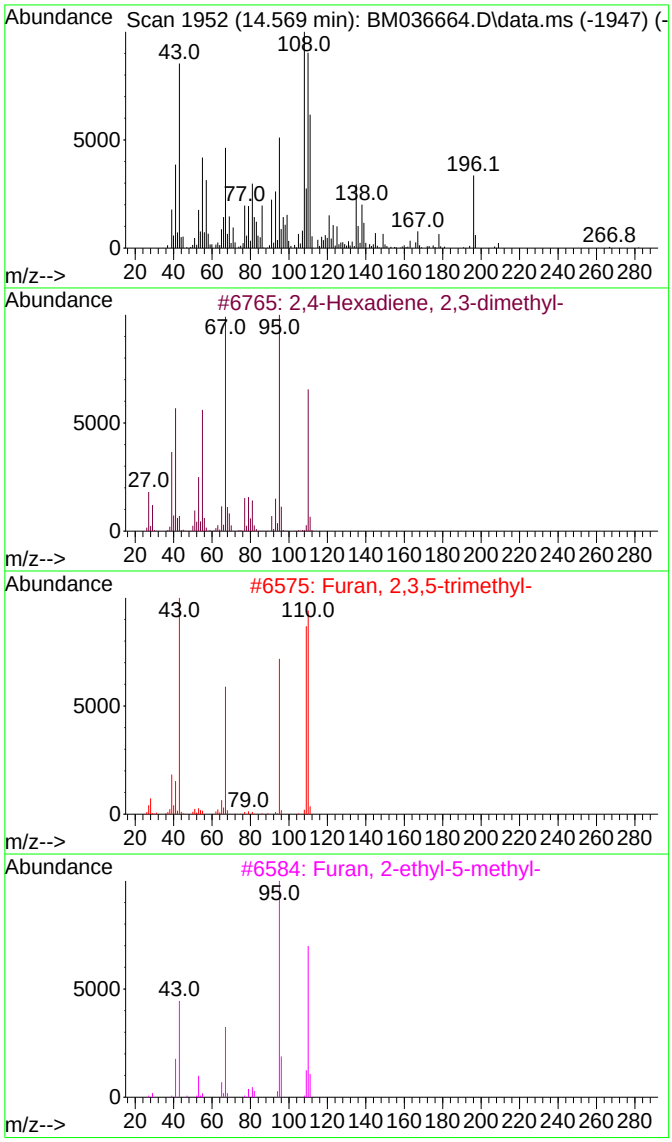
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

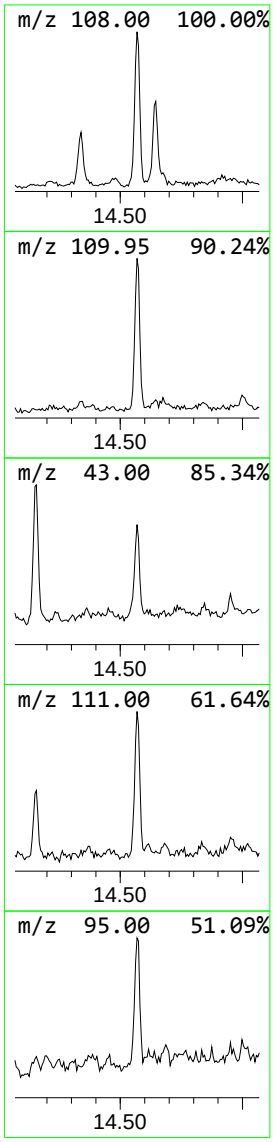
****

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.569	2.25 ng/ul	299252	Acenaphthene-d10	14.645

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Hexadiene, 2,3-dimethyl-	110	C8H14	005678-98-8	38
2	Furan, 2,3,5-trimethyl-	110	C7H10O	010504-04-8	30
3	Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	25
4	4-Ethyl-2-hexynal	124	C8H12O	071932-97-3	25
5	4,4-Dimethyl-2-cyclopenten-1-one	110	C7H10O	022748-16-9	22





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Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

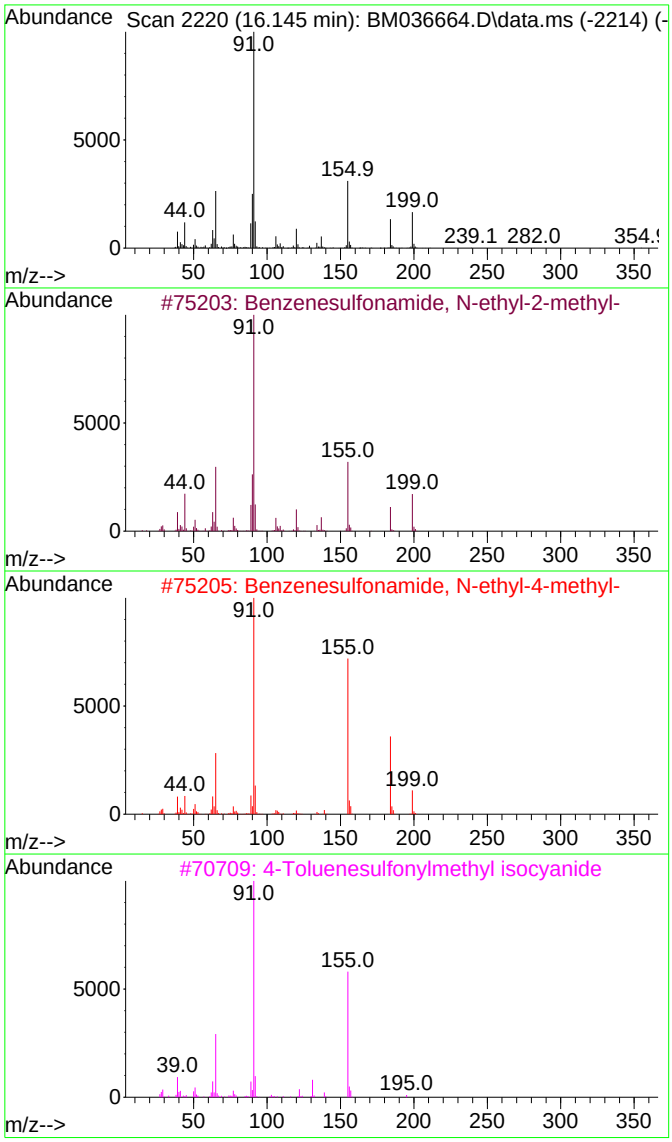
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

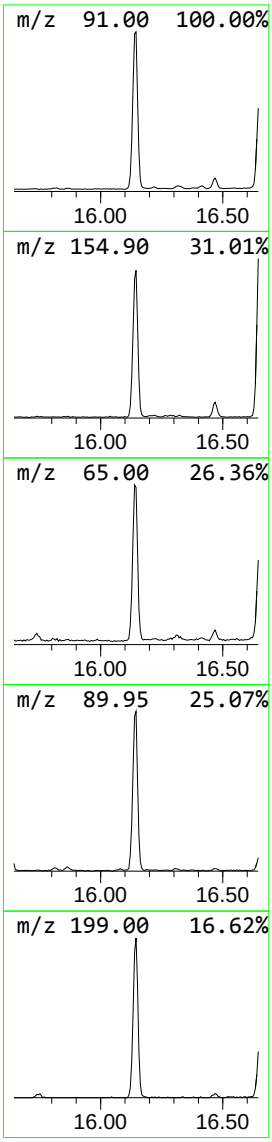
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	8.46 ng/ul	1203690	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	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	59
3	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50





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Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

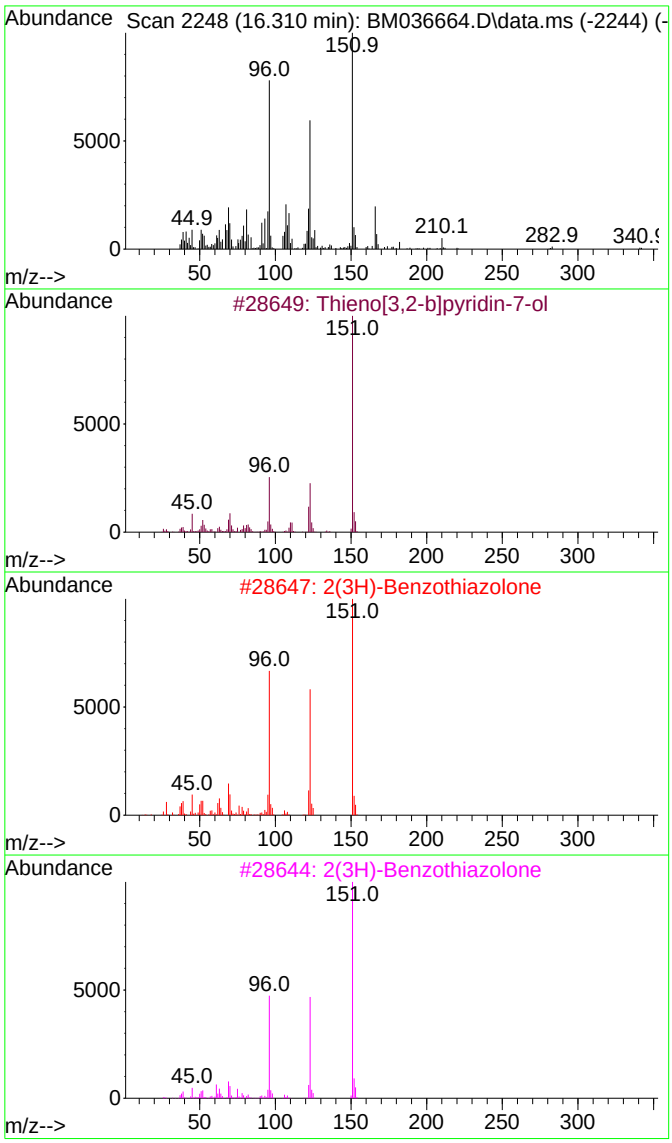
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

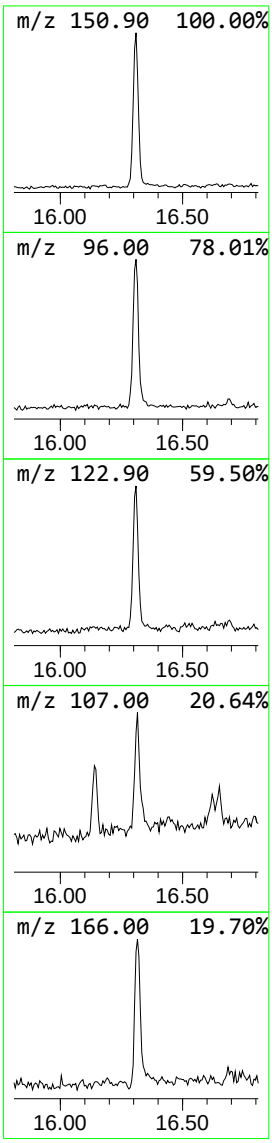
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Thieno[3,2-b]pyridin-7-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.310	2.36 ng/ul	336273	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	64
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	64
3	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	53
4	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	53
5	t-Butylhydroquinone	166	C10H14O2	001948-33-0	46





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Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

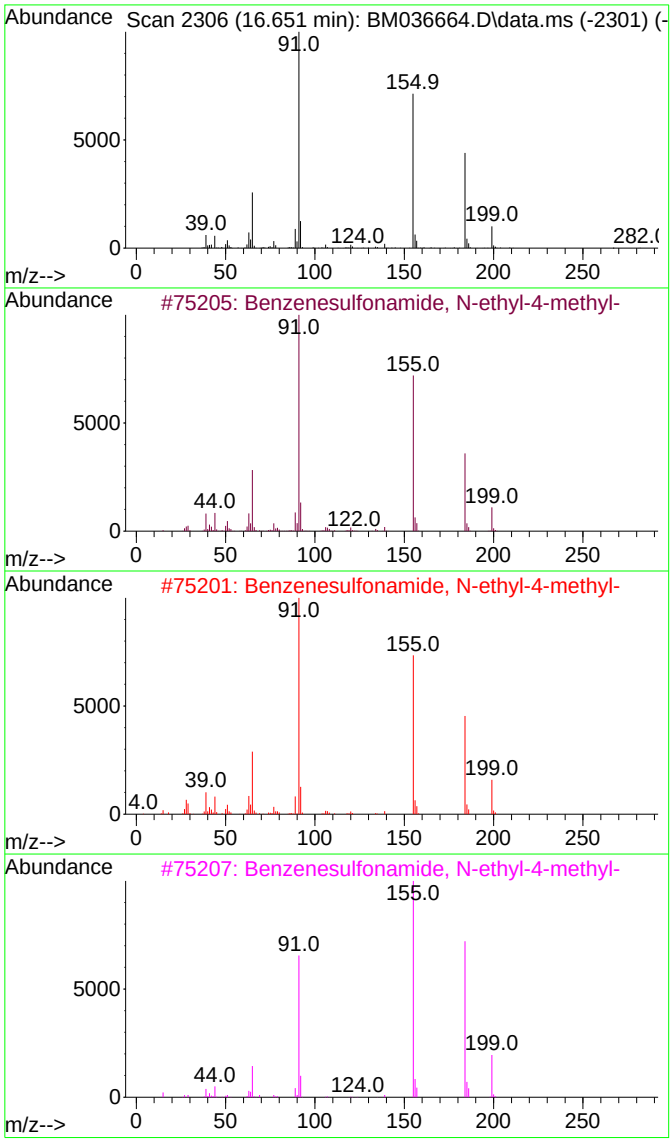
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

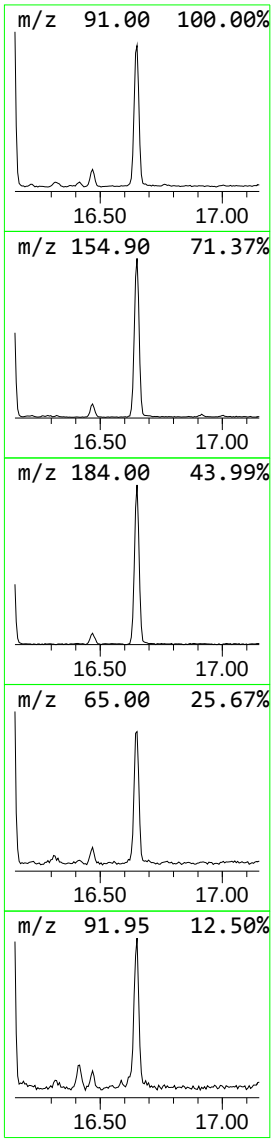
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	4.24 ng/ul	602560	Phenanthrene-d10	17.380

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	94
4	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
5	Benzene, 4-(ethylsulfonyl)-1-met...	184	C9H12O2S	007569-34-8	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

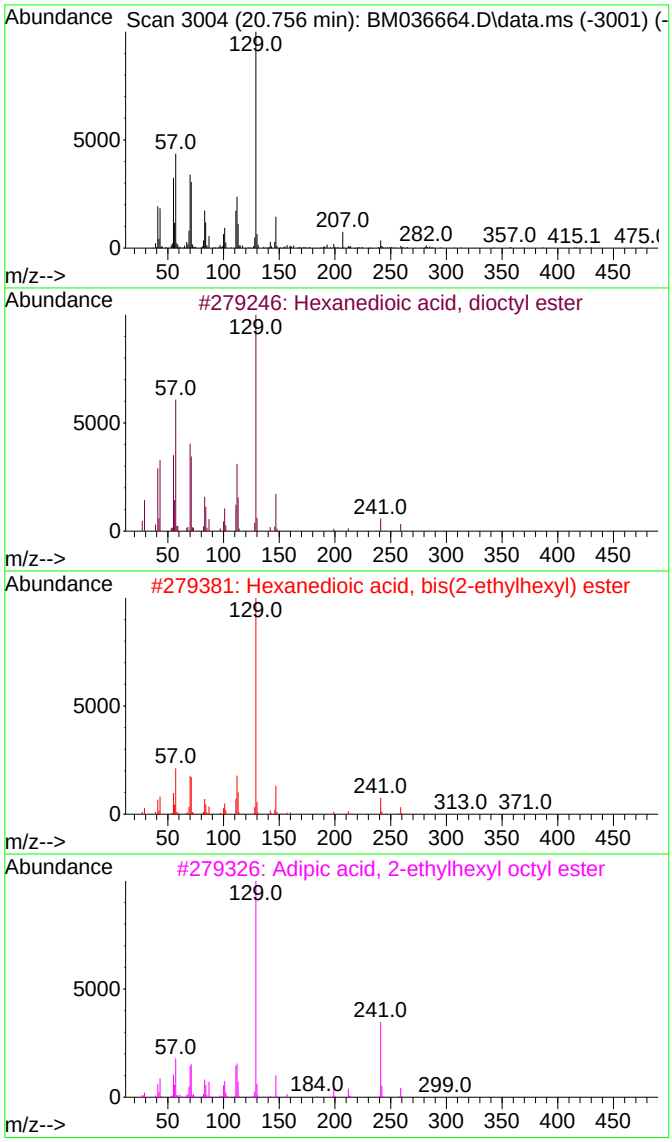
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

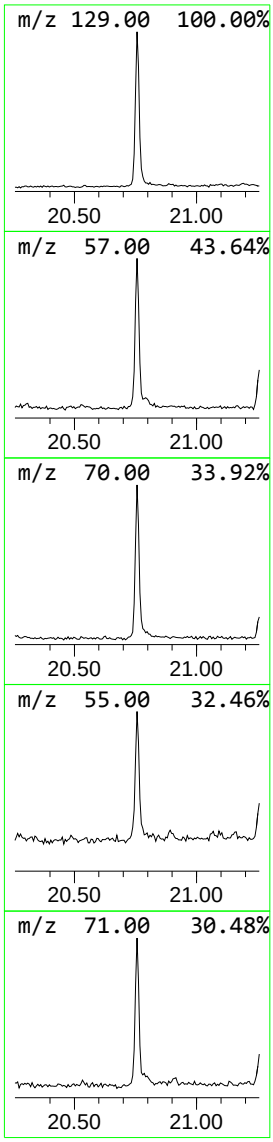
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Hexanedioic acid, dioctyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.756	3.28 ng/ul	358454	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	91
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
3	Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	90
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036664.D
Acq On : 22 Sep 2022 15:58
Operator : CG/JU
Sample : N4595-04DL 5X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8G9DL

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	31.5	ng/ul	2283670	1	8.016	1449730	20.0
Paraldehyde	4.234	2.3	ng/ul	169628	1	8.016	1449730	20.0
1,3-Oxathiolane	4.610	2.7	ng/ul	196730	1	8.016	1449730	20.0
2H-Pyran-3(4H)-...	6.716	2.1	ng/ul	151694	1	8.016	1449730	20.0
Benzenamine, 3-...	12.328	2.9	ng/ul	305086	2	10.834	2136640	20.0
unknown-01	14.569	2.3	ng/ul	299252	3	14.645	2660640	20.0
Benzenesulfonam...	16.145	8.5	ng/ul	1203690	4	17.380	2844740	20.0
Thieno[3,2-b]py...	16.310	2.4	ng/ul	336273	4	17.380	2844740	20.0
Benzenesulfonam...	16.651	4.2	ng/ul	602560	4	17.380	2844740	20.0
Hexanedioic aci...	20.756	3.3	ng/ul	358454	5	21.545	2183130	20.0