LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8K3DL

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: BM036748.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	13	rVB	3522623	3845165	79.66%	5.384%
2	3.181	13	16	27	rVB	135725	203888	4.22%	0.285%
3	3.422	53	57	65	rVB	393054	525995	10.90%	0.736%
4	3.669	96	99	106	rVB2	28531	40105	0.83%	0.056%
5	3.816	121	124	137	rBV	349343	546777	11.33%	0.766%
6	4.052	159	164	169	rBV	21542	35765	0.74%	0.050%
7	4.146	176	180	185	rVB	28333	40423	0.84%	0.057%
8	4.599	251	257	268	rVB	186561	307427	6.37%	0.430%
9	5.293	369	375	383	rVB	287118	426681	8.84%	0.597%
10	5.563	417	421	428	rVB5	16613	34135	0.71%	0.048%
11	5.734	446	450	458	rVV	49846	72170	1.50%	0.101%
12	5.928	478	483	488	rVV	17404	29907	0.62%	0.042%
13	6.269	536	541	546	rVB	20958	34861	0.72%	0.049%
14	6.481	568	577	583	rBV	36263	60840	1.26%	0.085%
15	6.587	590	595	600	rBV5	23582	45984	0.95%	0.064%
16	7.163	686	693	706	rVB	171902	333175	6.90%	0.466%
17	7.334	715	722	730	rBV	660264	1036156	21.47%	1.451%
18	7.534	745	756	765	rVB	470050	765088	15.85%	1.071%
19	7.663	774	778	785	rVB7	19276	35081	0.73%	0.049%
20	8.004	829	836	845	rVV	1400869	2198775	45.55%	3.078%
21	8.087	845	850	861	rVB5	43776	116420	2.41%	0.163%
22	8.381	893	900	907	rVB	299261	482095	9.99%	0.675%
23	8.551	923	929	933	rVB7	13024	26958	0.56%	0.038%
24	8.663	940	948	950	rBV4	28021	46420	0.96%	0.065%
25	8.710	950	956	965	rVV	497707	823833	17.07%	1.153%
26	8.992	998	1004	1013	rBV	443346	716697	14.85%	1.003%
27	9.116	1019	1025	1028	rBV2	71478	113408	2.35%	0.159%
28	9.169	1028	1034	1044	rVB	760438	1241683	25.73%	1.738%
29	9.375	1063	1069	1076	rVB10	32287	71294	1.48%	0.100%
30	9.528	1089	1095	1101	rBV2	67829	137356	2.85%	0.192%
31	9.634	1111	1113	1120	rVB	30712	45238	0.94%	0.063%
32	9.763	1130	1135	1139	rVB6	26139	40761	0.84%	0.057%
33	9.892	1151	1157	1170	rVB	348600	605190	12.54%	0.847%
34	10.004	1170	1176	1177	rBV4	25184	36379	0.75%	0.051%
35	10.034	1177	1181	1188	rVV6	33138	72536	1.50%	0.102%
36	10.210	1203	1211	1222	rBV6	46792	119195	2.47%	0.167%

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
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Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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	10.363	1230	1237	1242	rVB5	85020	217297	4.50%	0.304%
38	10.434	1242	1249	1259	rBV	699502	1197224	24.80%	1.676%
39	10.604	1271	1278	1287	rBV8	25766	67266	1.39%	0.094%
40	10.739	1296	1301	1307	rBV	120141	210330	4.36%	0.294%
41	10.816	1307	1314	1321	rVV	1921318	3205357	66.41%	4.488%
42	10.951	1329	1337	1352	rVV	924697	1637334	33.92%	2.292%
43	11.069	1352	1357	1361	rVV4	37649	73700	1.53%	0.103%
44	11.116	1361	1365	1372	rVV4	43358	80943	1.68%	0.113%
45	11.181	1372	1376	1379	rVB5	18780	28107	0.58%	0.039%
46	11.398	1403	1413	1419	rVB5	20909	60270	1.25%	0.084%
47	11.604	1439	1448	1451	rBV5	31503	75901	1.57%	0.106%
48	11.786	1475	1479	1484	rVB5	26413	45411	0.94%	0.064%
49	12.016	1515	1518	1523	rVB2	28164	42197	0.87%	0.059%
50	12.151	1535	1541	1545	rBV	164446	254014	5.26%	0.356%
51	12.310	1563	1568	1573	rBV2	133196	210801	4.37%	0.295%
52	12.369	1574	1578	1580	rVV4	26073	42313	0.88%	0.059%
53	12.410	1581	1585	1598	rVB4	45801	152059	3.15%	0.213%
54	12.522	1599	1604	1610	rBV6	30232	63622	1.32%	0.089%
55	12.686	1628	1632	1636	rVB4	46883	68835	1.43%	0.096%
56	12.798	1646	1651	1656	rBV2	50662	81327	1.68%	0.114%
57	12.898	1664	1668	1676	rVB10	33445	69125	1.43%	0.097%
58	13.022	1679	1689	1692	rBV10	30515	92801	1.92%	0.130%
59	13.369	1744	1748	1753	rVB8	22195	38601	0.80%	0.054%
60	13.533	1772	1776	1782	rVB	101434	148830	3.08%	0.208%
61	13.692	1798	1803	1807	rBV4	33378	60092	1.24%	0.084%
62	13.774	1813	1817	1819	rBV4	21310	37262	0.77%	0.052%
63	13.957	1843	1848	1852	rBV4	54053	116392	2.41%	0.163%
64	14.039	1857	1862	1872	rVB	1838156	2528663	52.39%	3.540%
65	14.186	1883	1887	1893	rBV6	54952	117336	2.43%	0.164%
66	14.321	1905	1910	1924	rVB	1793992	2782613	57.65%	3.896%
67	14.433	1925	1929	1936	rVB5	37394	70711	1.46%	0.099%
68	14.557	1945	1950	1955	rVB	258437	342615	7.10%	0.480%
69	14.627	1956	1962	1968	rVV	2935923	4279524	88.66%	5.992%
70	14.692	1968	1973	1980	rVV	507606	756466	15.67%	1.059%
71	14.757	1981	1984	1991	rVV	52211	83483	1.73%	0.117%
72	14.827	1992	1996	2003	rVB	91495	139299	2.89%	0.195%
73	15.027	2025	2030	2038	rVB	257496	399572	8.28%	0.559%
74	15.368	2083	2088	2093	rBV	263598	389760	8.08%	0.546%
75	15.616	2124	2130	2135	rBV	2583788	3642932	75.47%	5.100%
76	15.668	2135	2139	2145	rVB	373343	525518	10.89%	0.736%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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

77	15.727	2145	2149	2154	rBV	256958	346061	7.17%	0.485%
78	15.845	2165	2169	2172	rBV3	48245	93822	1.94%	0.131%
79	16.127	2211	2217	2230	rVB	1180433	1695990	35.14%	2.375%
80	16.298	2240	2246	2252	rBV2	249460	399024	8.27%	0.559%
81	16.457	2270	2273	2277	rVB	48723	57272	1.19%	0.080%
82	16.639	2298	2304	2308	rBV	619677	884550	18.33%	1.238%
83	16.904	2346	2349	2352	rBV	62780	70960	1.47%	0.099%
84	17.174	2391	2395	2401	rVB2	99953	182688	3.78%	0.256%
85	17.368	2422	2428	2432	rBV	3451726	4826715	100.00%	6.758%
86	17.410	2432	2435	2439	rVV	480032	619969	12.84%	0.868%
87	17.462	2439	2444	2455	rVB	2779853	3961366	82.07%	5.546%
88	18.039	2538	2542	2546	rVB	128215	159268	3.30%	0.223%
89	18.262	2575	2580	2590	rBV2	403051	699018	14.48%	0.979%
90	19.033	2708	2711	2721	rVB	141697	201281	4.17%	0.282%
91	19.415	2773	2776	2781	rVB	172045	212520	4.40%	0.298%
92	19.468	2783	2785	2789	rVB2	76236	82373	1.71%	0.115%
93	19.533	2792	2796	2801	rBV	163404	234195	4.85%	0.328%
94	19.751	2827	2833	2837	rBV	3133450	4269923	88.46%	5.978%
95	19.792	2837	2840	2848	rVB2	449640	632950	13.11%	0.886%
96	20.745	2999	3002	3004	rBV	198591	217354	4.50%	0.304%
97	21.245	3084	3087	3095	rBV	865949	1161841	24.07%	1.627%
98	21.533	3131	3136	3143	rBV2	3254150	4269779	88.46%	5.978%
99	23.809	3517	3523	3530	rBV2	1430831	2810660	58.23%	3.935%
100	23.968	3545	3550	3563	rVB2	1773619	3585529	74.29%	5.020%

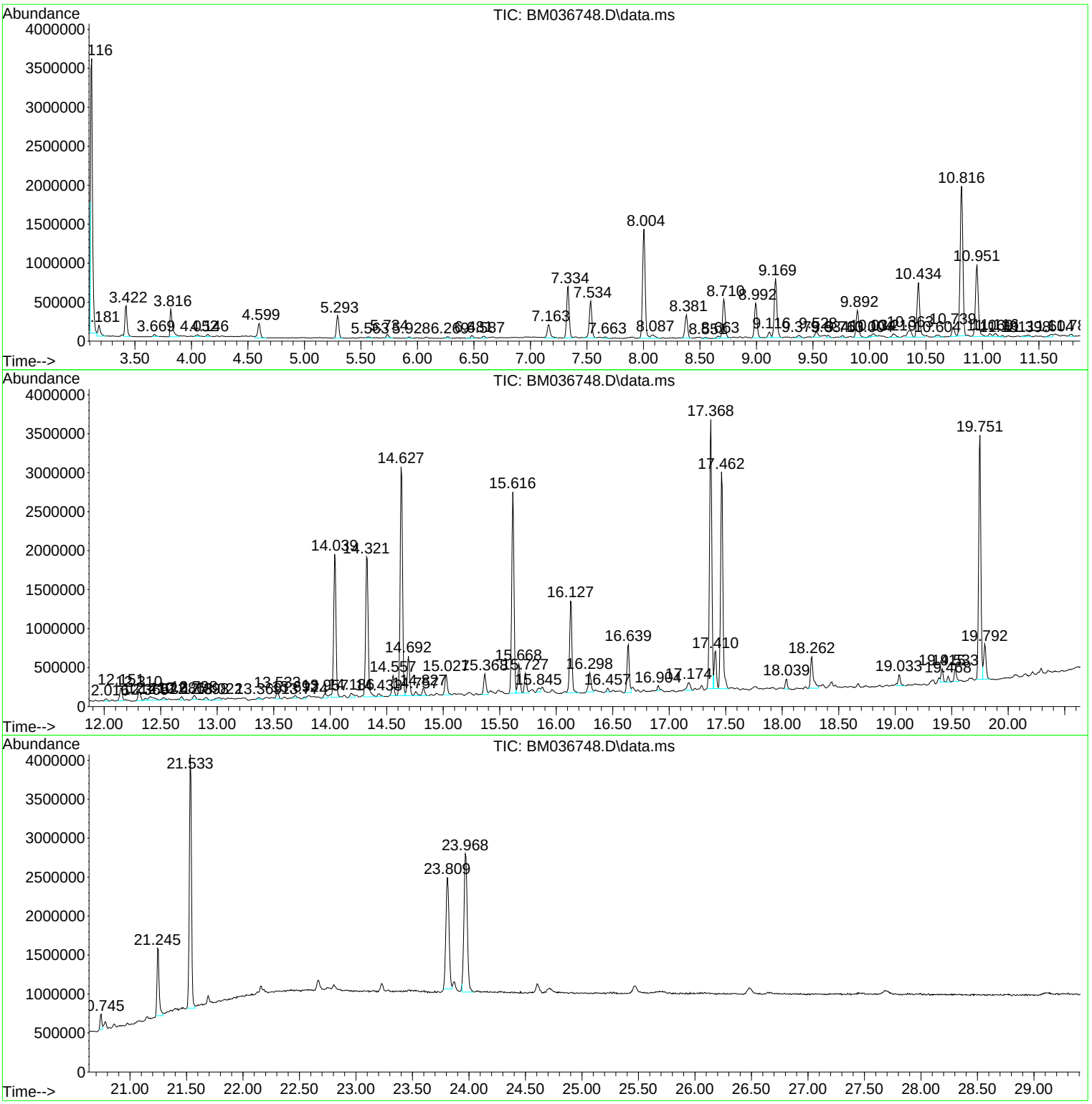
Sum of corrected areas: 71424872

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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

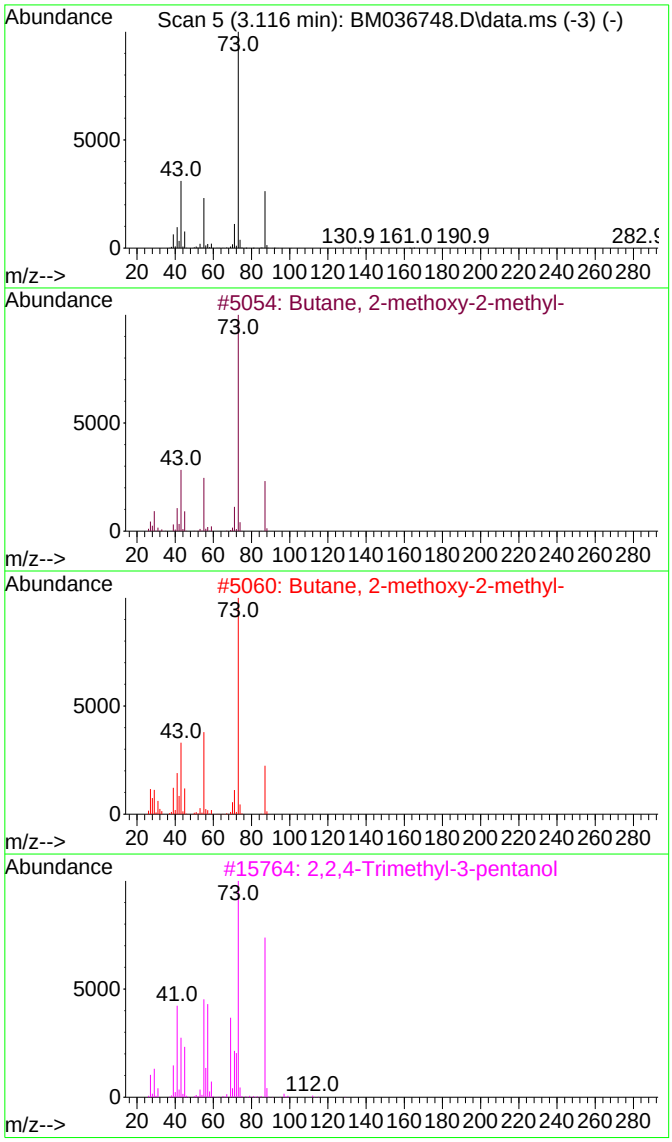
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

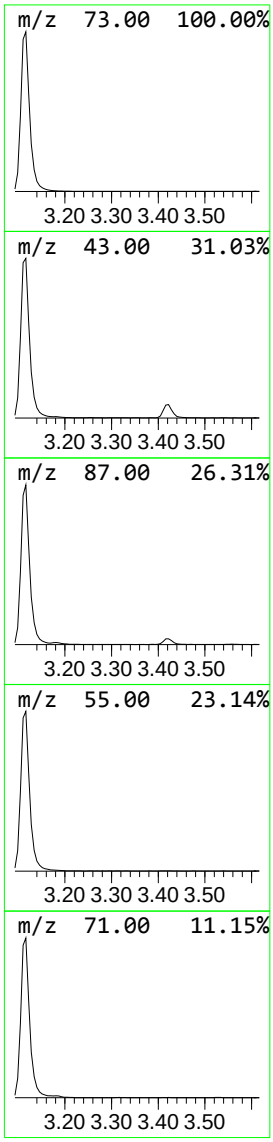
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.116	34.98 ng/ul	3845170	1,4-Dichlorobenzene-d4	8.004

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	40
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

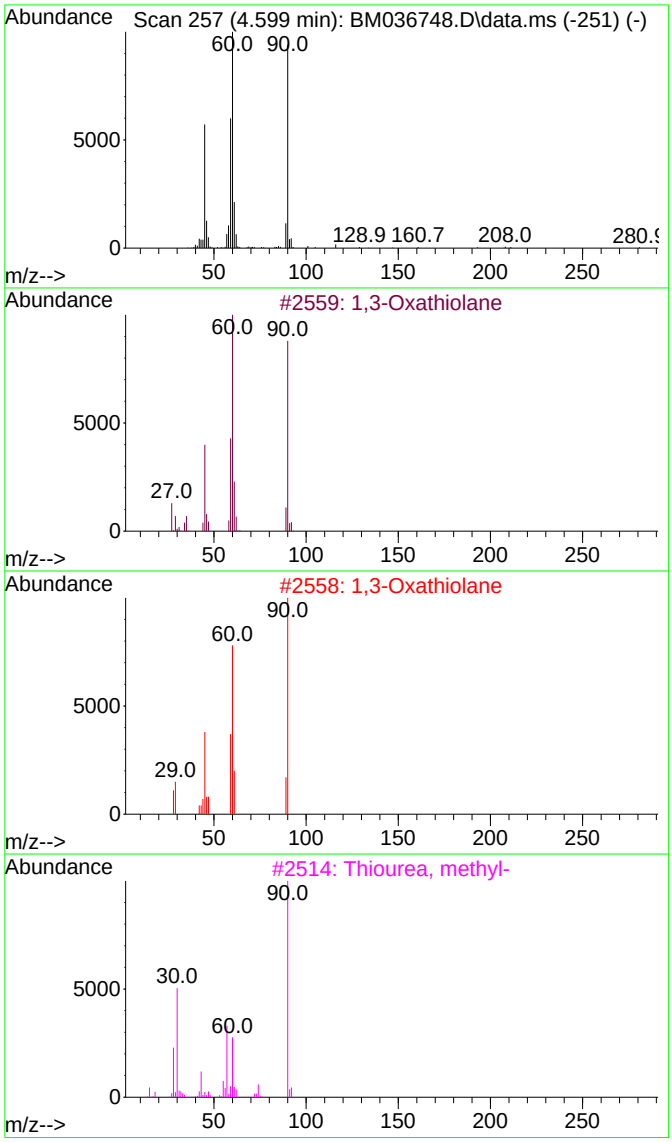
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

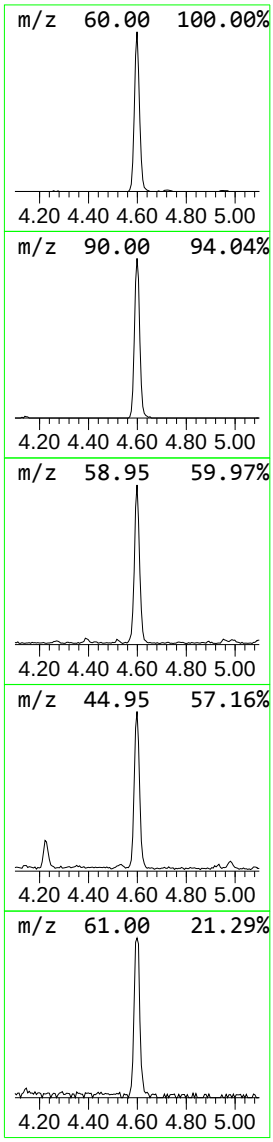
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1,3-Oxathiolane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.599	2.80 ng/ul	307427	1,4-Dichlorobenzene-d4	8.004

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	91
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	56
3	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
4	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	45
5	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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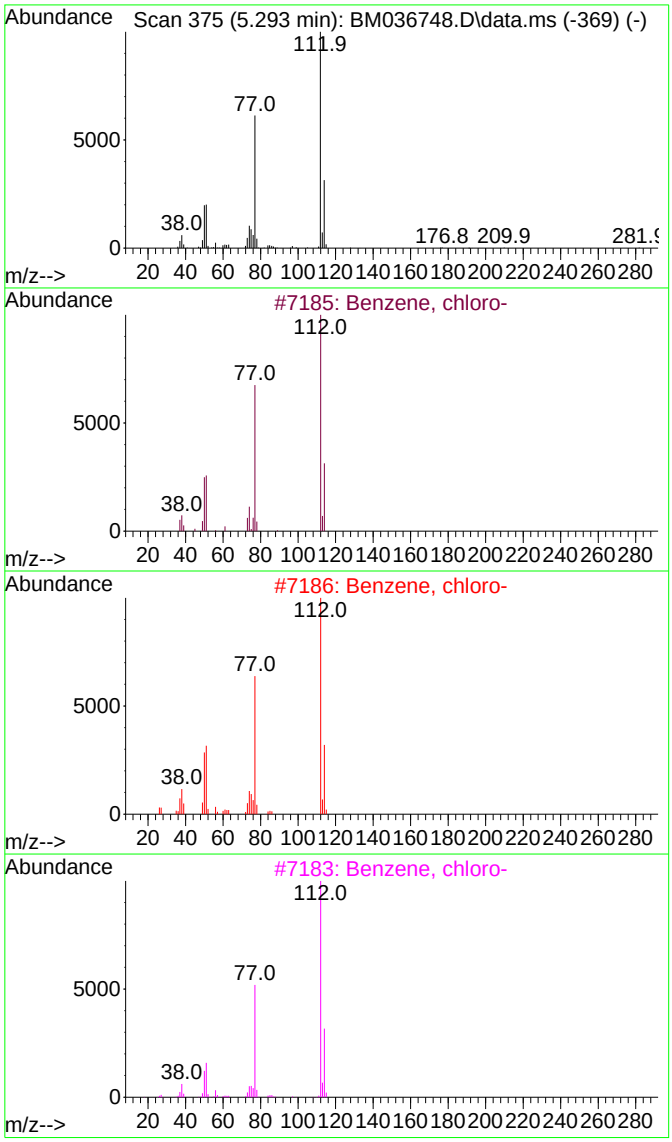
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

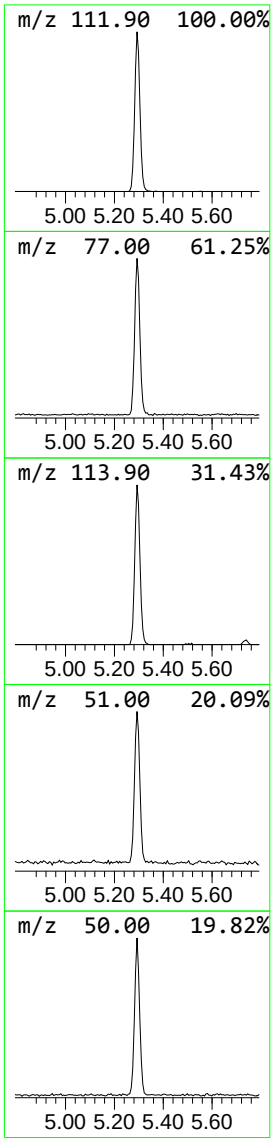
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.293	3.88 ng/ul	426681	1,4-Dichlorobenzene-d4	8.004

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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

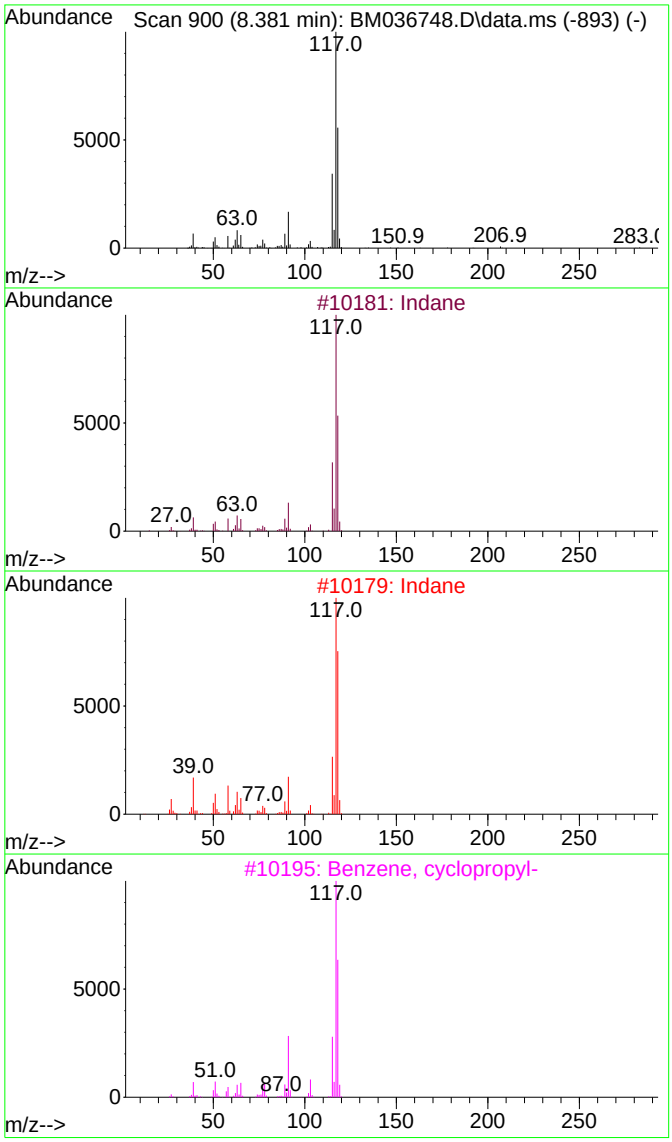
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

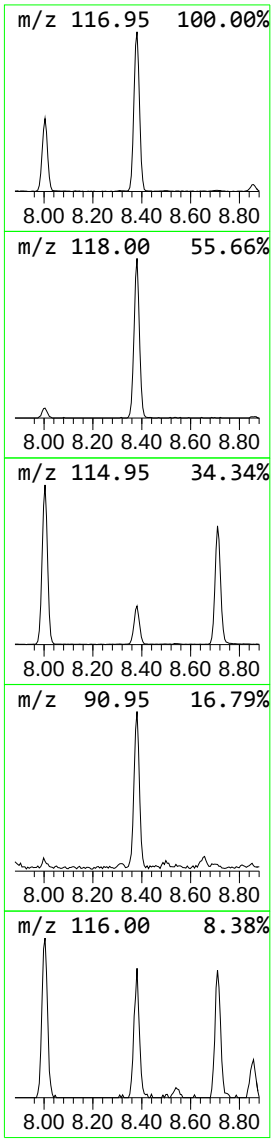
****

Peak Number 4 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.381	4.39 ng/ul	482095	1,4-Dichlorobenzene-d4	8.004

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

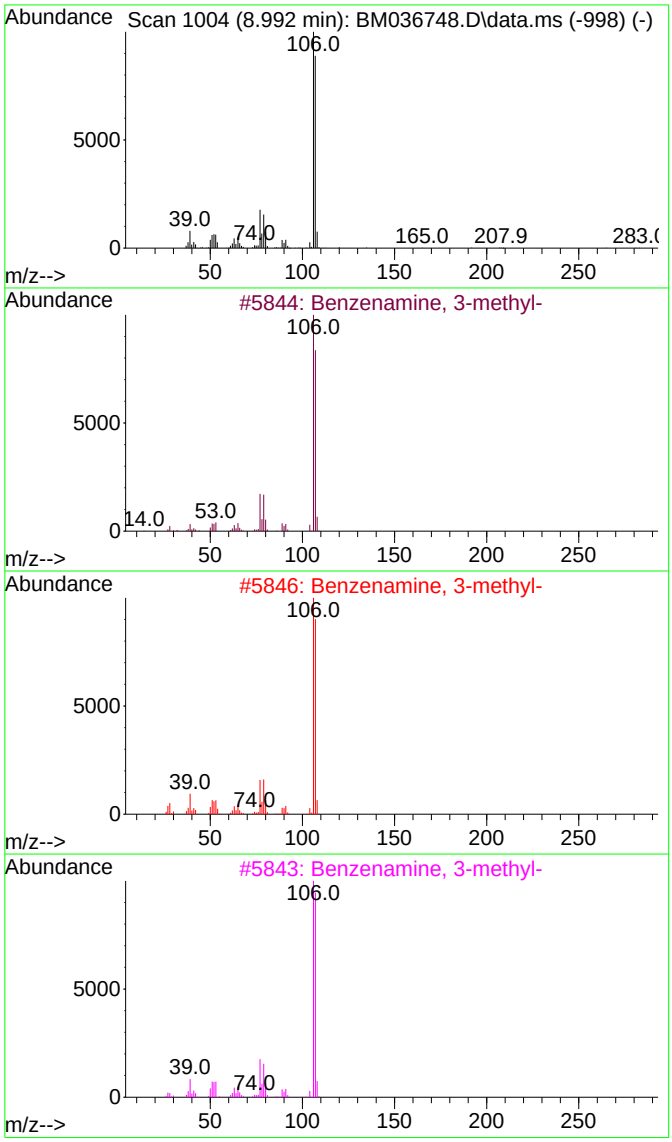
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

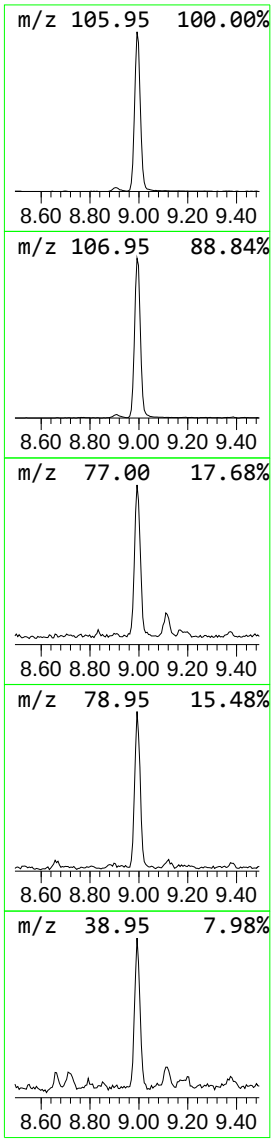
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzenamine, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.992	6.52 ng/ul	716697	1,4-Dichlorobenzene-d4	8.004

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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Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

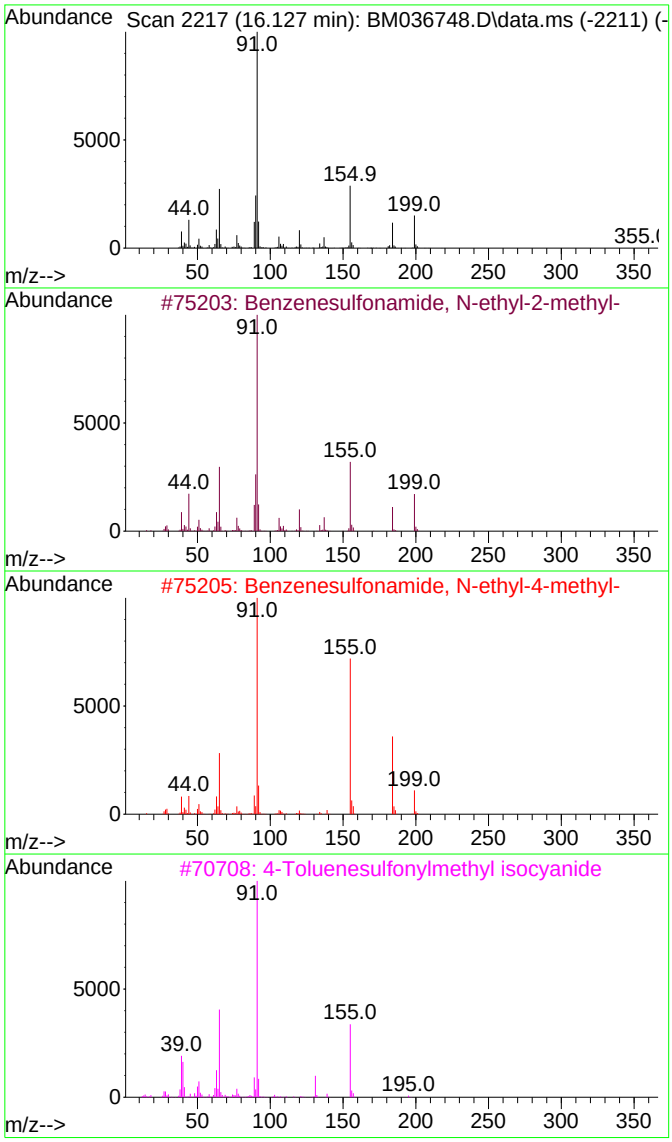
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

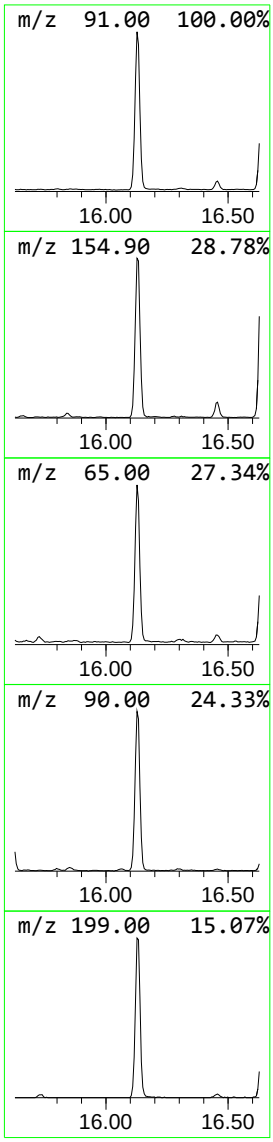
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.127	7.03 ng/ul	1695990	Phenanthrene-d10	17.368

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

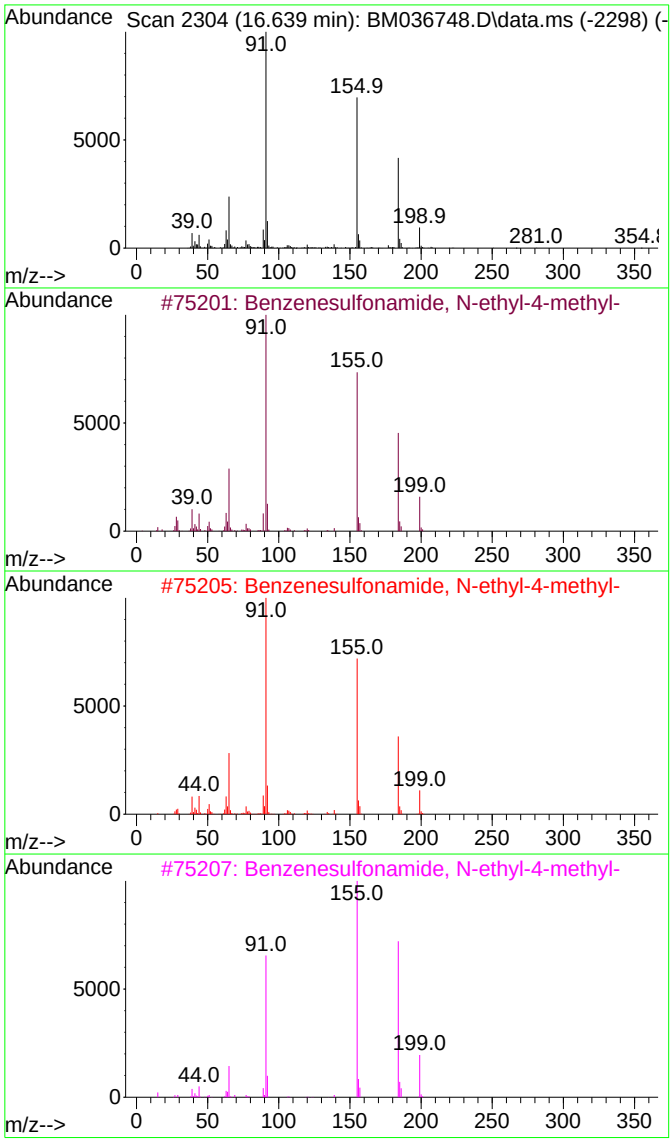
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

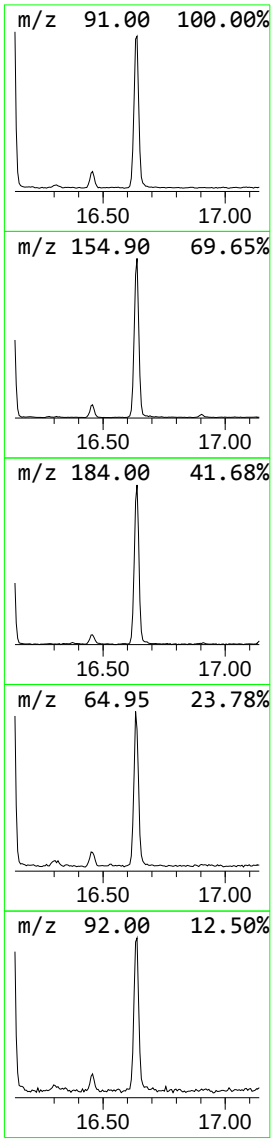
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	3.67 ng/ul	884550	Phenanthrene-d10	17.368

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	97
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	94
4	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
5	Benzene, 4-(ethylsulfonyl)-1-met...	184	C9H12O2S	007569-34-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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

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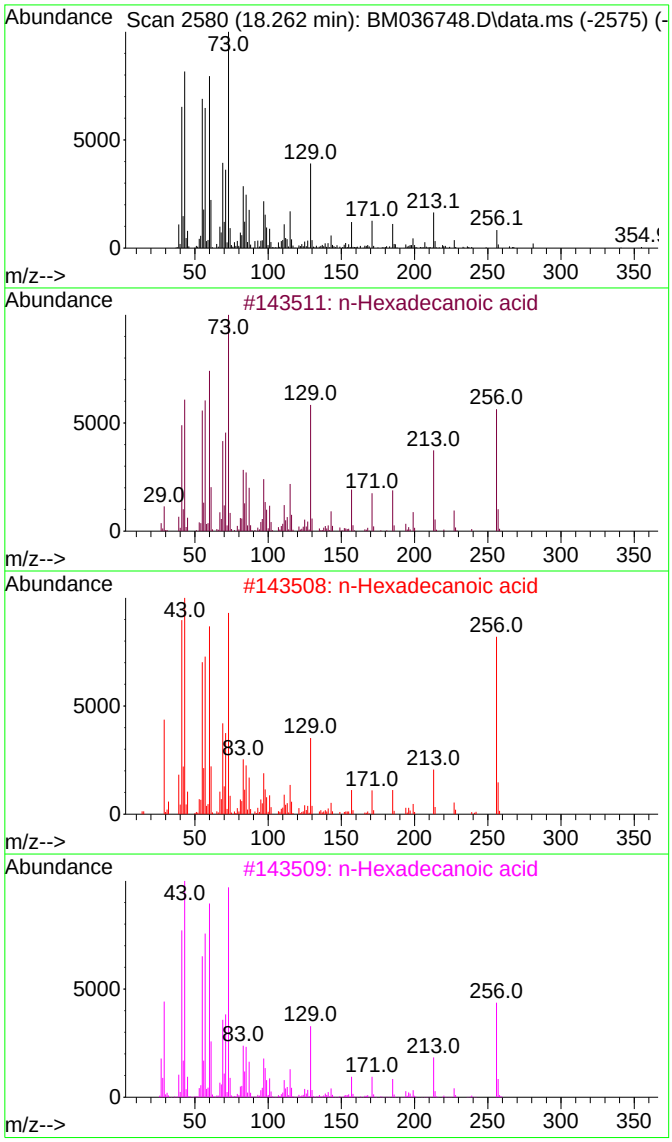
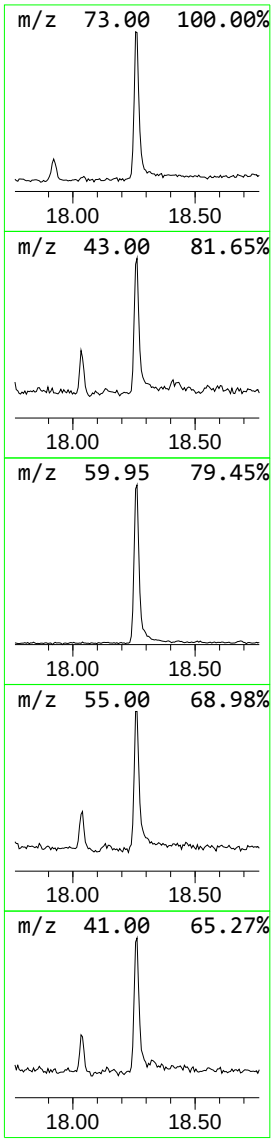
Peak Number 8 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.262	2.90 ng/ul	699018	Phenanthrene-d10	17.368

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	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91





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Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

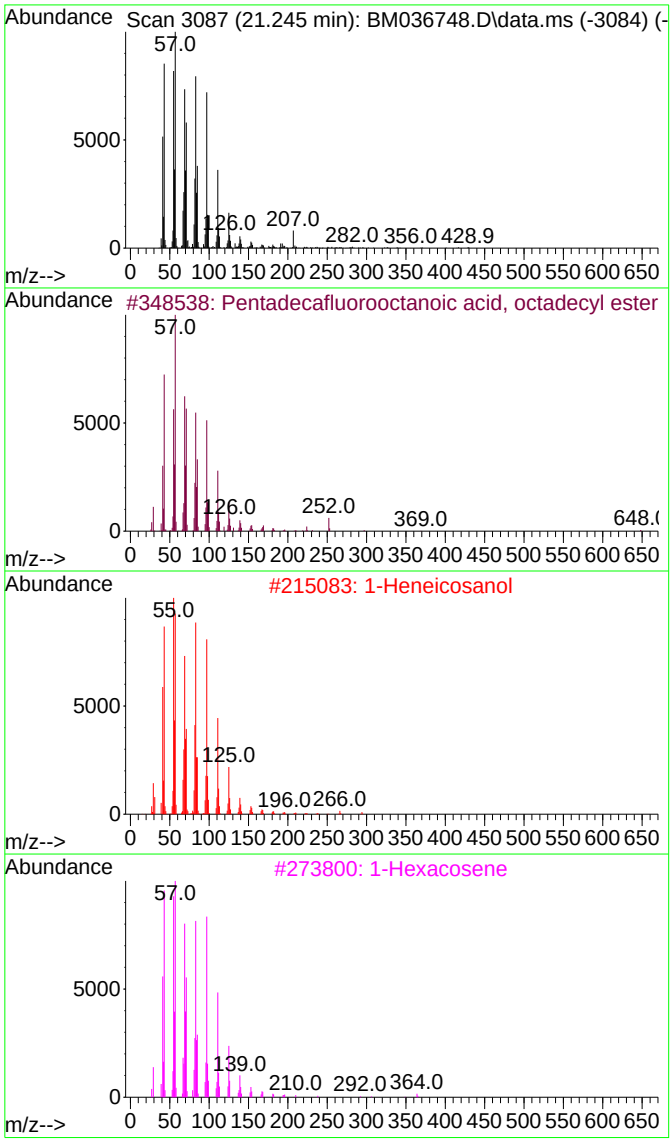
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

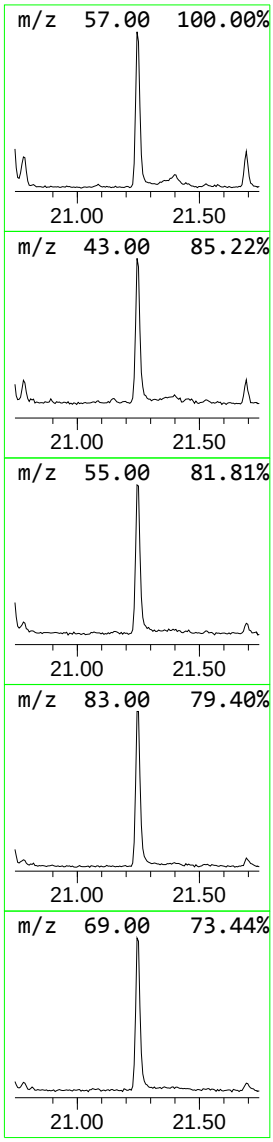
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Pentadecafluorooctanoic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.245	5.44 ng/ul	1161840	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	1-Hexacosene	364	C26H52	018835-33-1	91
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5	Nonadecyl heptafluorobutyrate	480	C23H39F7O2	1000351-83-9	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
Data File : BM036748.D
Acq On : 26 Sep 2022 16:49
Operator : CG/JU
Sample : N4706-03DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8K3DL

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...	3.116	35.0	ng/ul	3845170	1	8.004	2198780	20.0
1,3-Oxathiolane	4.599	2.8	ng/ul	307427	1	8.004	2198780	20.0
Benzene, chloro-	5.293	3.9	ng/ul	426681	1	8.004	2198780	20.0
Indane	8.381	4.4	ng/ul	482095	1	8.004	2198780	20.0
Benzenamine, 3-...	8.992	6.5	ng/ul	716697	1	8.004	2198780	20.0
Benzenesulfonam...	16.127	7.0	ng/ul	1695990	4	17.368	4826720	20.0
Benzenesulfonam...	16.639	3.7	ng/ul	884550	4	17.368	4826720	20.0
n-Hexadecanoic ...	18.262	2.9	ng/ul	699018	4	17.368	4826720	20.0
Pentadecafluoro...	21.245	5.4	ng/ul	1161840	5	21.533	4269780	20.0