LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033978.D
Acq On : 02 Feb 2022 17:50
Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

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-BM013122.M
Title : SVOA CALIBRATION

Signal : TIC: BM033978.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.284	5	8	11	rBV	22000	27736	1.06%	0.081%
2	3.307	11	12	18	rVB2	16449	18925	0.73%	0.055%
3	3.584	55	59	64	rVV6	10402	14846	0.57%	0.043%
4	3.643	64	69	80	rVV	942345	1231480	47.23%	3.575%
5	3.719	80	82	100	rVV	63348	116101	4.45%	0.337%
6	3.872	104	108	114	rVV	82496	111439	4.27%	0.323%
7	3.954	114	122	130	rVV	63474	128811	4.94%	0.374%
8	4.054	134	139	150	rVB	81264	118005	4.53%	0.343%
9	4.419	198	201	206	rVB3	14283	19659	0.75%	0.057%
10	4.466	206	209	217	rBV	17051	25311	0.97%	0.073%
11	5.001	293	300	311	rVV	388489	543669	20.85%	1.578%
12	5.348	353	359	363	rVV7	8434	20662	0.79%	0.060%
13	5.407	363	369	378	rVV	193624	280565	10.76%	0.814%
14	5.543	386	392	406	rVB	613195	1194881	45.83%	3.469%
15	5.748	423	427	437	rVB	34948	54623	2.10%	0.159%
16	5.919	450	456	463	rVV	316087	484943	18.60%	1.408%
17	5.990	464	468	478	rVB	32564	47502	1.82%	0.138%
18	6.190	496	502	506	rBV7	7854	16912	0.65%	0.049%
19	6.401	530	538	545	rBV7	8388	18323	0.70%	0.053%
20	6.560	561	565	569	rBV5	9481	16603	0.64%	0.048%
21	6.913	615	625	631	rBV6	15124	35760	1.37%	0.104%
22	6.972	631	635	638	rVV6	5528	9953	0.38%	0.029%
23	7.007	638	641	647	rVV	22302	34542	1.32%	0.100%
24	7.078	647	653	662	rVV	435070	746412	28.63%	2.167%
25	7.142	662	664	672	rVB	14306	19604	0.75%	0.057%
26	7.248	675	682	689	rBV	339707	526883	20.21%	1.529%
27	7.307	689	692	697	rVB	9912	13345	0.51%	0.039%
28	7.454	707	717	725	rBV	431767	710695	27.26%	2.063%
29	7.525	725	729	730	rVV3	10054	12582	0.48%	0.037%
30	7.566	730	736	743	rVB3	31160	74937	2.87%	0.218%
31	7.713	757	761	766	rVV2	11615	18290	0.70%	0.053%
32	7.925	789	797	804	rBV	415496	662286	25.40%	1.923%
33	7.989	804	808	814	rBV2	8178	12203	0.47%	0.035%
34	8.042	814	817	824	rVB3	10026	13785	0.53%	0.040%
35	8.160	827	837	845	rBV	63499	114941	4.41%	0.334%
36	8.472	886	890	895	rVB5	5054	9828	0.38%	0.029%

LSC Area Percent Report

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Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

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-BM013122.M
Title : SVOA CALIBRATION

37	8.631	904	917	926	rVV	560758	921841	35.36%	2.676%
38	8.707	926	930	932	rVV5	5540	10205	0.39%	0.030%
39	8.748	932	937	944	rVB	27717	51433	1.97%	0.149%
40	9.036	978	986	988	rBV6	7288	13326	0.51%	0.039%
41	9.084	988	994	1001	rVV	435089	714022	27.39%	2.073%
42	9.166	1005	1008	1012	rVV	15083	22416	0.86%	0.065%
43	9.213	1012	1016	1021	rVV	16706	26314	1.01%	0.076%
44	9.536	1065	1071	1077	rVV	26515	41890	1.61%	0.122%
45	9.813	1110	1118	1126	rBV	359056	599523	22.99%	1.740%
46	9.972	1140	1145	1153	rVV2	17368	35302	1.35%	0.102%
47	10.266	1191	1195	1203	rVB8	6386	10847	0.42%	0.031%
48	10.354	1203	1210	1226	rBV	556153	952767	36.54%	2.766%
49	10.519	1233	1238	1246	rVB	19214	34405	1.32%	0.100%
50	10.736	1267	1275	1283	rBV	604540	1009002	38.70%	2.929%
51	10.866	1290	1297	1314	rBV	422857	760665	29.18%	2.208%
52	12.166	1514	1518	1522	rBV2	12757	20316	0.78%	0.059%
53	12.207	1522	1525	1530	rVV	10806	19726	0.76%	0.057%
54	12.260	1530	1534	1540	rVV	12664	22230	0.85%	0.065%
55	12.324	1540	1545	1550	rVB2	9683	17219	0.66%	0.050%
56	12.977	1652	1656	1661	rVB8	6285	9941	0.38%	0.029%
57	13.401	1722	1728	1734	rBV3	34877	50907	1.95%	0.148%
58	13.971	1819	1825	1831	rBV	1120894	1480364	56.78%	4.297%
59	14.254	1867	1873	1881	rVB	1130526	1669281	64.02%	4.846%
60	14.466	1903	1909	1915	rVB	23971	30925	1.19%	0.090%
61	14.560	1919	1925	1935	rVV	983609	1455686	55.83%	4.226%
62	14.748	1949	1957	1972	rBV	317616	593990	22.78%	1.724%
63	14.971	1990	1995	2001	rVB8	5797	10048	0.39%	0.029%
64	15.077	2007	2013	2020	rVB9	3681	10542	0.40%	0.031%
65	15.195	2028	2033	2040	rVB3	8737	13354	0.51%	0.039%
66	15.371	2056	2063	2066	rBV6	7896	16599	0.64%	0.048%
67	15.413	2066	2070	2074	rVB3	14562	18566	0.71%	0.054%
68	15.548	2087	2093	2106	rBV	1486266	2165096	83.04%	6.285%
69	15.660	2106	2112	2126	rBV	451821	638279	24.48%	1.853%
70	15.789	2130	2134	2137	rBV5	9151	12724	0.49%	0.037%
71	15.965	2160	2164	2167	rBV2	9068	11622	0.45%	0.034%
72	16.236	2205	2210	2214	rBV	52621	80148	3.07%	0.233%
73	16.295	2218	2220	2224	rVV4	10454	12118	0.46%	0.035%
74	16.383	2231	2235	2240	rBV	10660	15358	0.59%	0.045%
75	16.448	2240	2246	2251	rBV4	18498	32526	1.25%	0.094%
76	16.507	2251	2256	2259	rBV2	12424	15902	0.61%	0.046%

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Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

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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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Title : SVOA CALIBRATION

77	16.542	2260	2262	2268	rVB5	7924	11378	0.44%	0.033%
78	16.618	2268	2275	2278	rBV6	10679	25631	0.98%	0.074%
79	16.701	2286	2289	2297	rVB6	13070	21609	0.83%	0.063%
80	16.771	2297	2301	2305	rBV	13221	22052	0.85%	0.064%
81	16.842	2310	2313	2319	rVB4	10081	14007	0.54%	0.041%
82	17.059	2346	2350	2354	rBV	14327	22982	0.88%	0.067%
83	17.201	2369	2374	2376	rBV6	6530	11247	0.43%	0.033%
84	17.295	2384	2390	2400	rBV	1226657	1790284	68.67%	5.197%
85	17.395	2400	2407	2419	rVV2	1645632	2367113	90.79%	6.871%
86	17.483	2419	2422	2428	rVB6	8493	16182	0.62%	0.047%
87	17.971	2502	2505	2509	rVB2	20736	23797	0.91%	0.069%
88	18.077	2519	2523	2526	rBV5	5463	10399	0.40%	0.030%
89	18.189	2538	2542	2549	rBV2	71417	111431	4.27%	0.323%
90	18.501	2590	2595	2602	rBV8	11632	24604	0.94%	0.071%
91	18.618	2612	2615	2620	rVB7	10205	15219	0.58%	0.044%
92	19.248	2718	2722	2726	rBV3	22749	35940	1.38%	0.104%
93	19.318	2731	2734	2737	rVV	18078	22345	0.86%	0.065%
94	19.465	2756	2759	2764	rBV5	21716	29530	1.13%	0.086%
95	19.683	2790	2796	2801	rBV	1949369	2607242	100.00%	7.568%
96	21.171	3045	3049	3057	rBV	144212	245039	9.40%	0.711%
97	21.459	3093	3098	3111	rBV	1365992	1819845	69.80%	5.283%
98	23.371	3418	3423	3429	rVB	515456	823712	31.59%	2.391%
99	23.694	3472	3478	3485	rBV2	941166	1810856	69.45%	5.257%
100	23.847	3498	3504	3518	rVB2	703579	1432126	54.93%	4.157%

Sum of corrected areas: 34449037

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033978.D
Acq On : 02 Feb 2022 17:50
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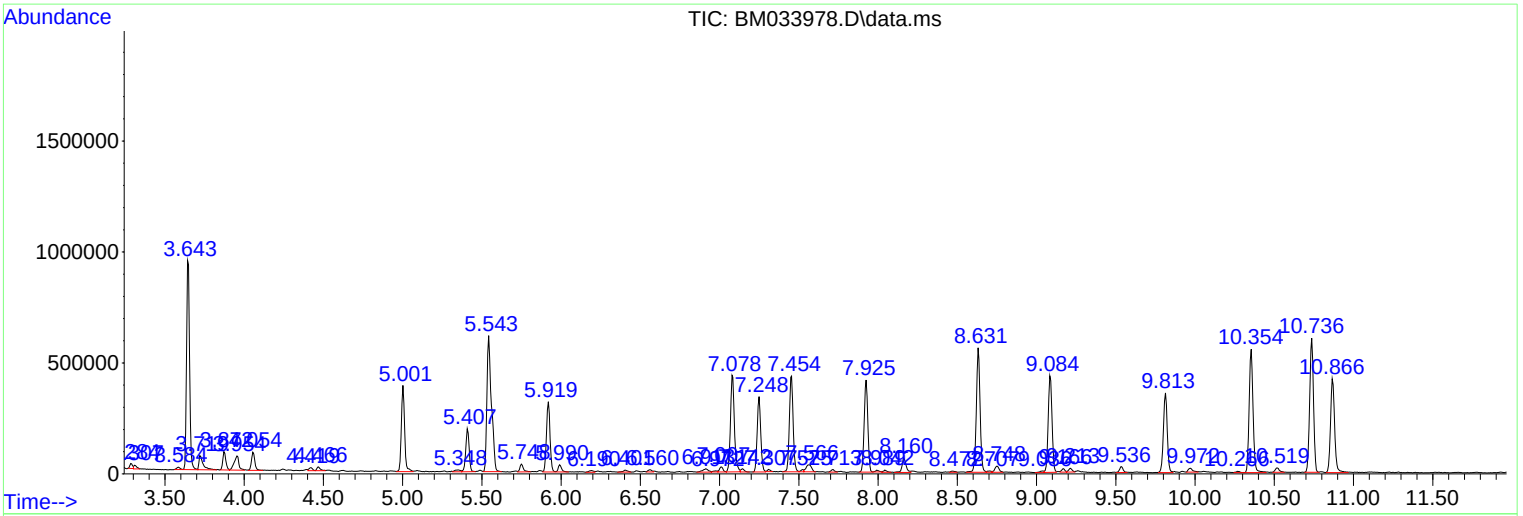
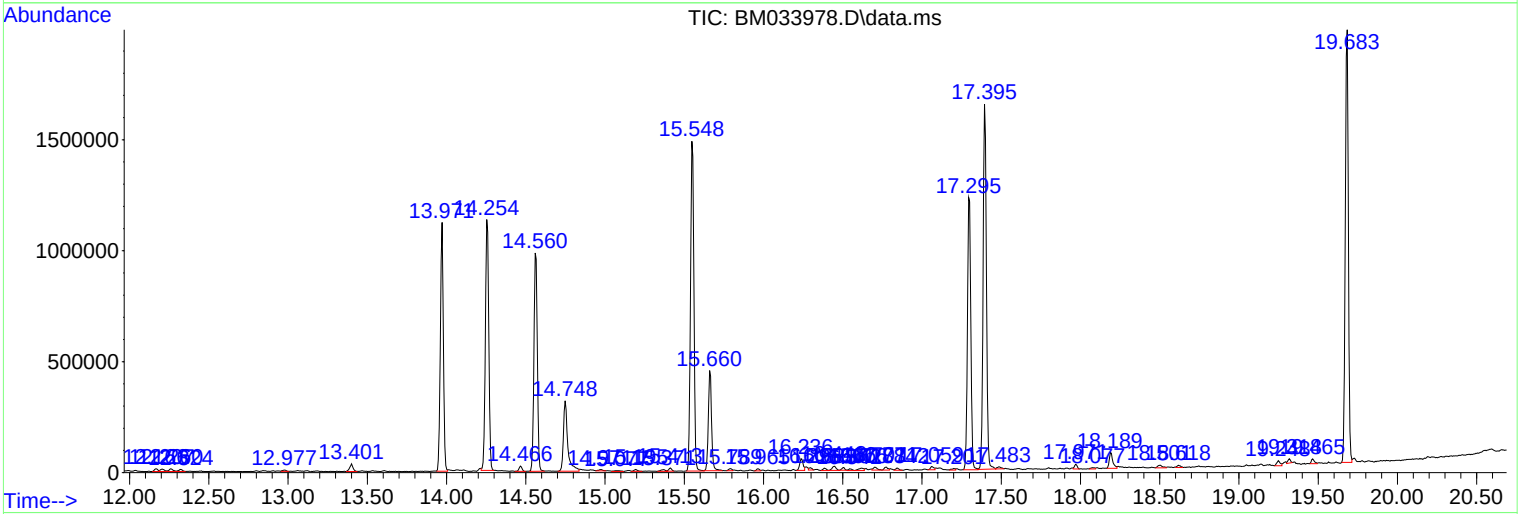
Misc :
ALS Vial : 49 Sample Multiplier: 1

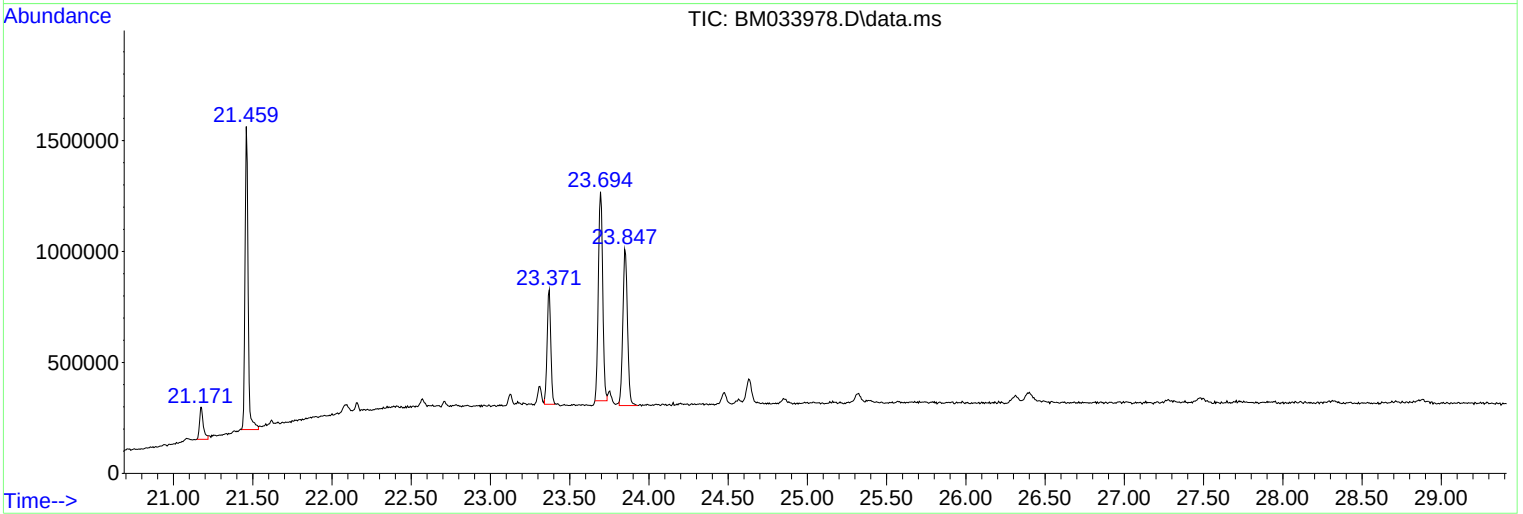
Instrument :
BNA_M
ClientSampleId :
C0VE2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033978.D
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Instrument :
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ClientSampleId :
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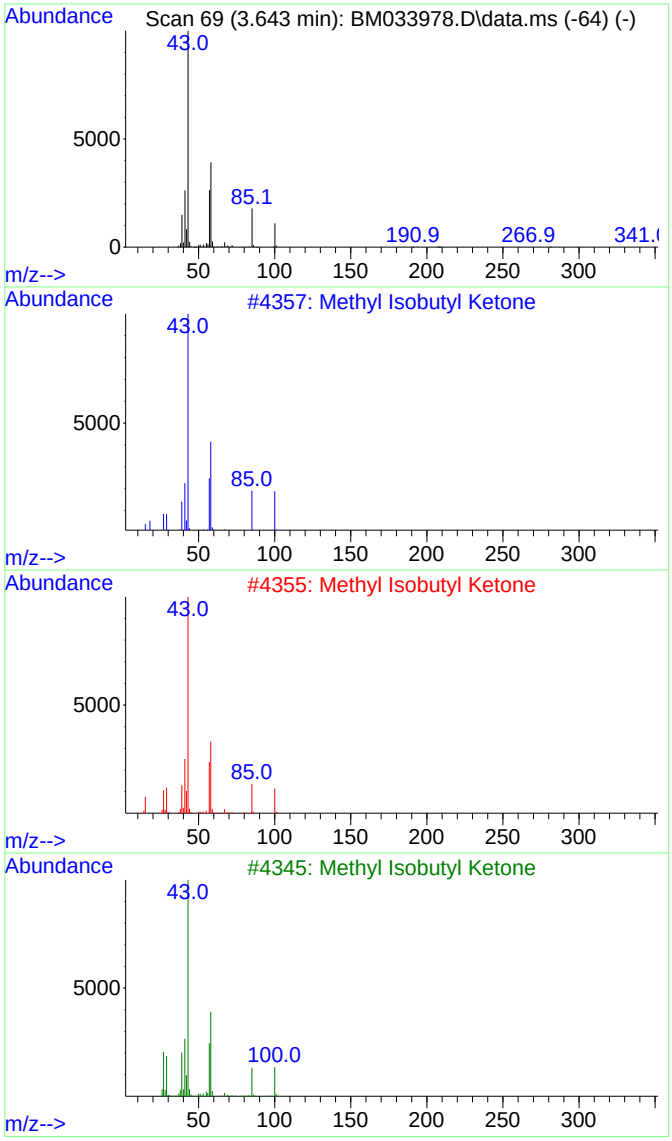
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

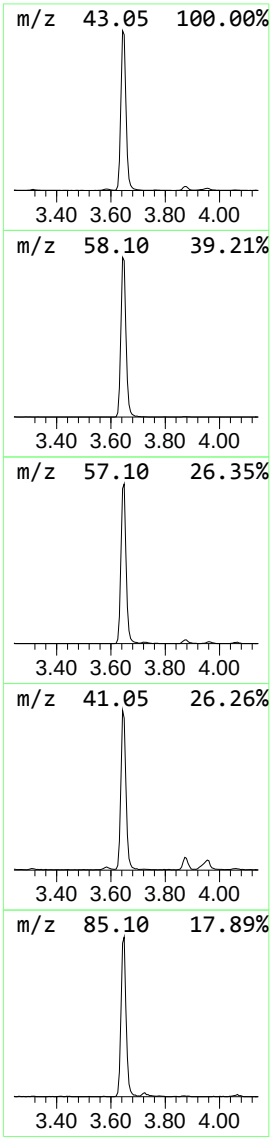
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.643	37.19 ng/ul	1231480	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	91
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	87
3	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	87
4	2-Hexanone	100	C6H12O	000591-78-6	78
5	2-Hexanone	100	C6H12O	000591-78-6	78





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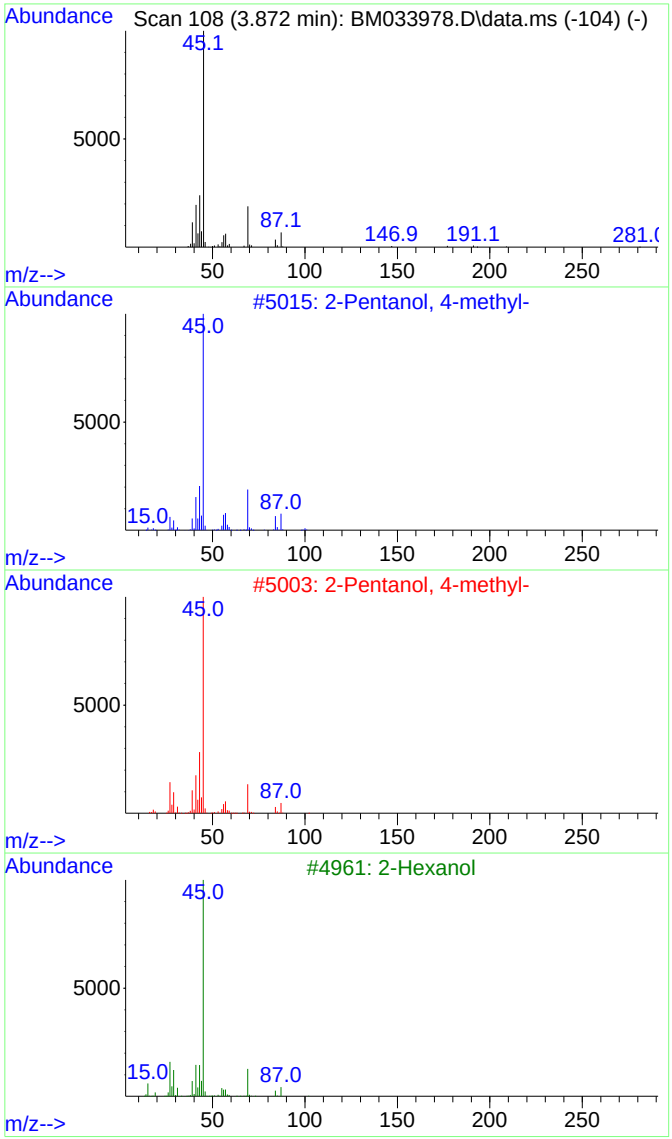
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

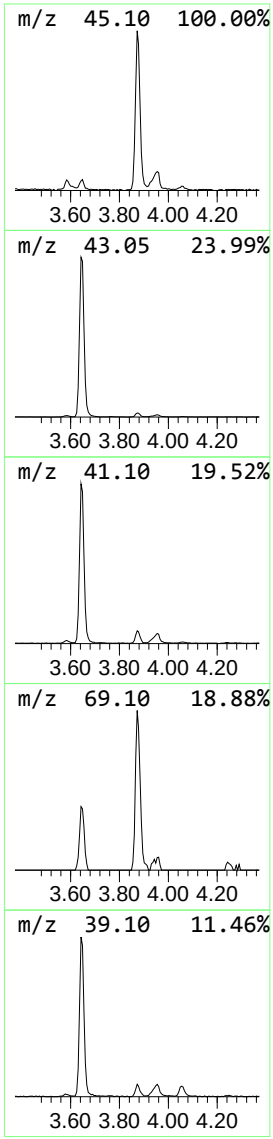
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanol, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.872	3.37 ng/ul	111439	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
3	2-Hexanol	102	C6H14O	000626-93-7	83
4	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
5	2-Octanol, (R)-	130	C8H18O	005978-70-1	78





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Sample : N1355-01
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ALS Vial : 49 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

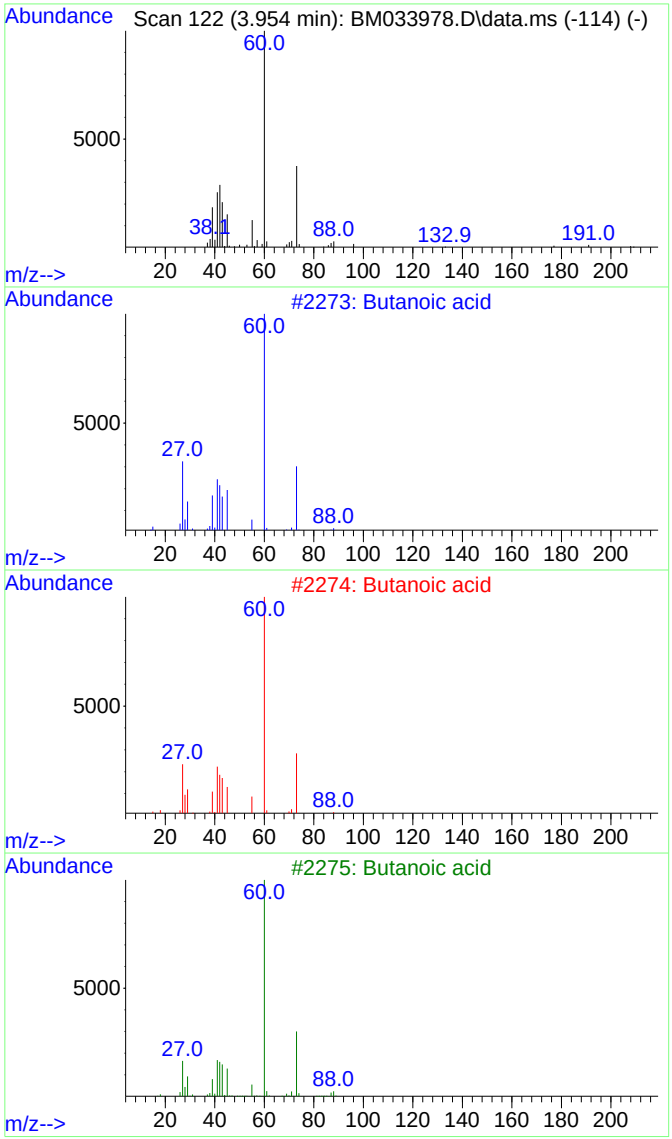
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

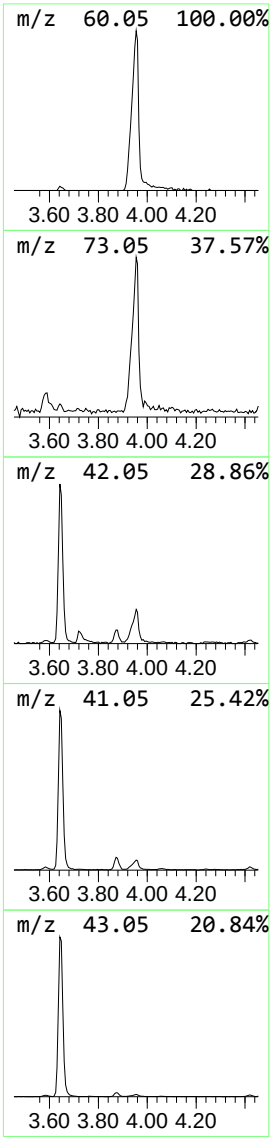
****

Peak Number 3 Butanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.954	3.89 ng/ul	128811	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	87
2	Butanoic acid	88	C4H8O2	000107-92-6	83
3	Butanoic acid	88	C4H8O2	000107-92-6	80
4	Butanoic acid	88	C4H8O2	000107-92-6	80
5	Pentanoic acid	102	C5H10O2	000109-52-4	56





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Data File : BM033978.D
Acq On : 02 Feb 2022 17:50
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Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
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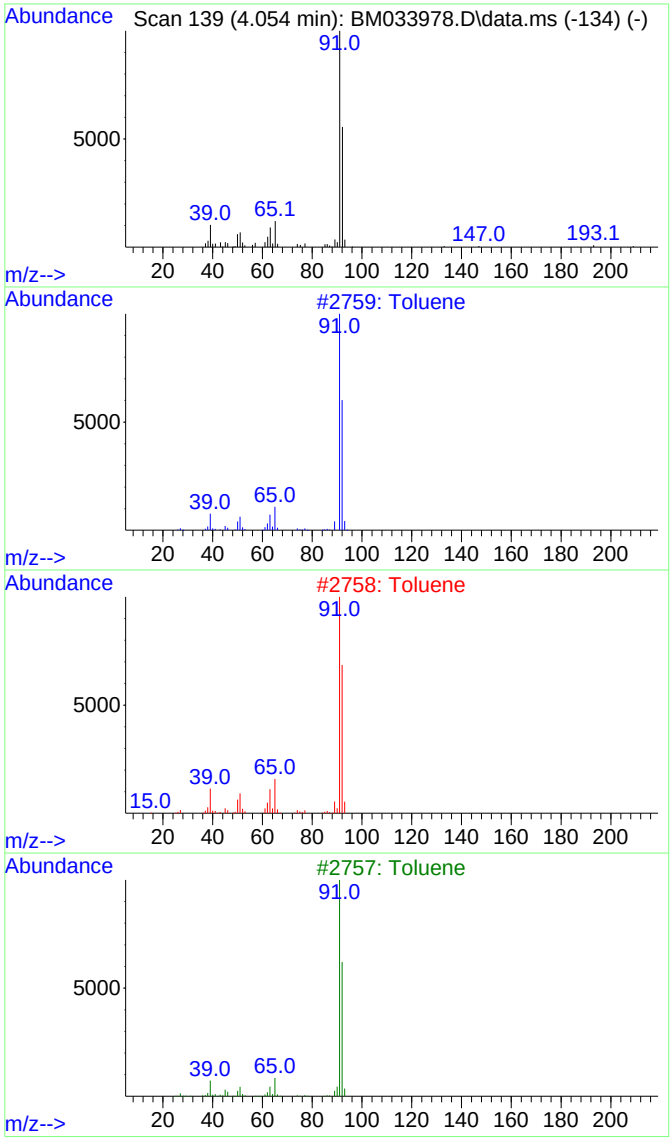
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

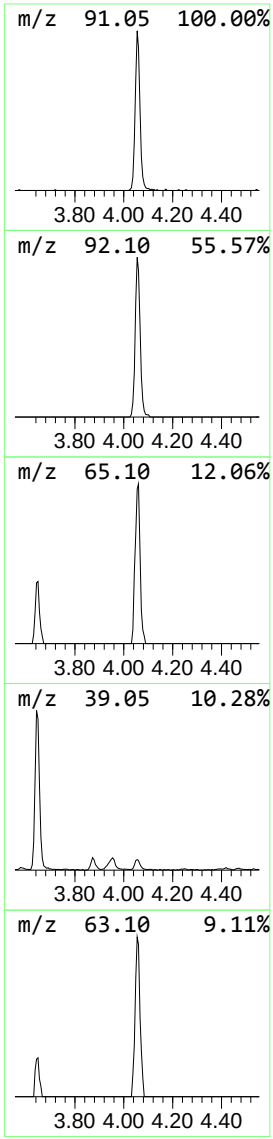
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Toluene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.054	3.56 ng/ul	118005	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	94
2	Toluene	92	C7H8	000108-88-3	93
3	Toluene	92	C7H8	000108-88-3	87
4	Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	87
5	Toluene	92	C7H8	000108-88-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033978.D
Acq On : 02 Feb 2022 17:50
Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

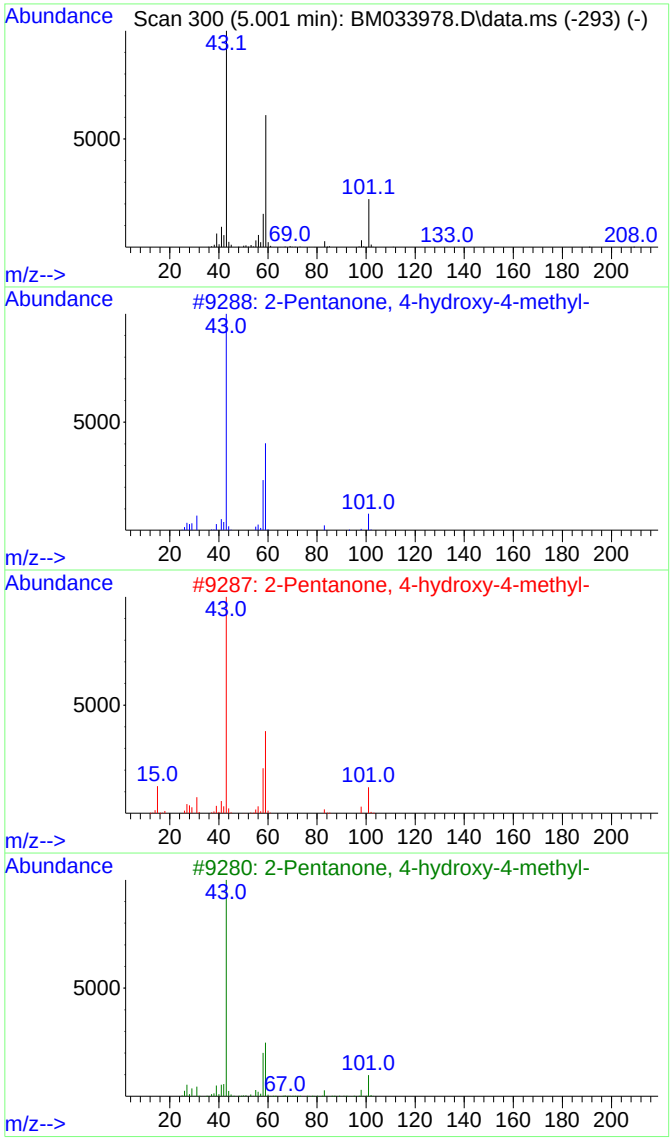
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

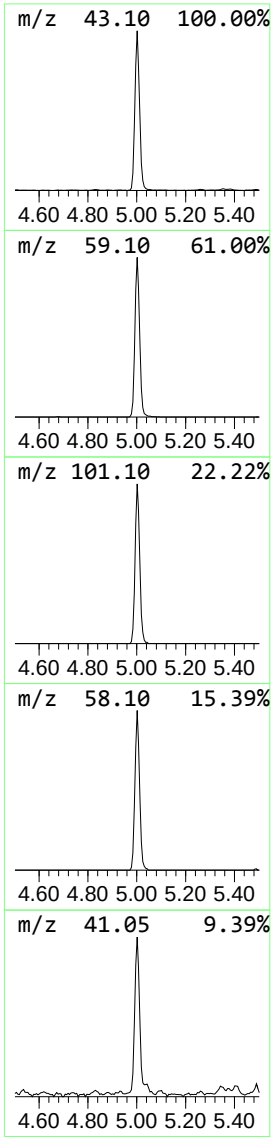
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.001	16.42 ng/ul	543669	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32





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Acq On : 02 Feb 2022 17:50
Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

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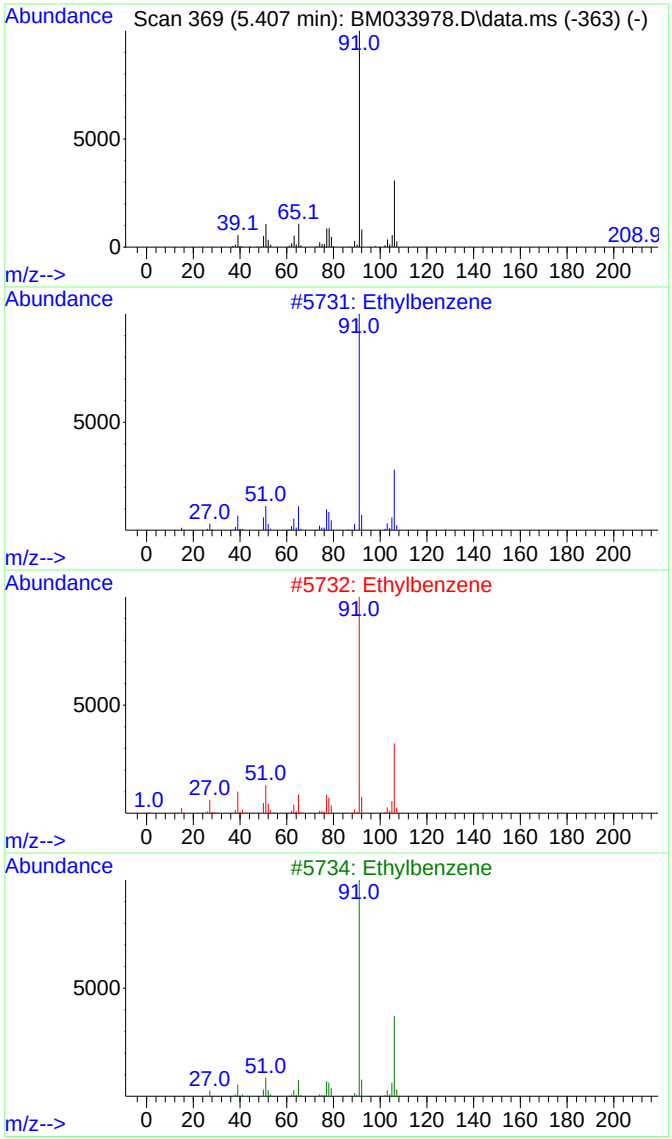
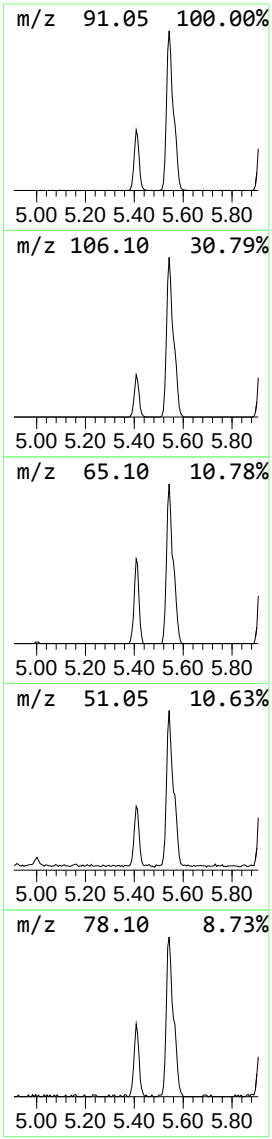
Peak Number 6 Ethylbenzene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.407	8.47 ng/ul	280565	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	95
2	Ethylbenzene	106	C8H10	000100-41-4	94
3	Ethylbenzene	106	C8H10	000100-41-4	93
4	Ethylbenzene	106	C8H10	000100-41-4	91
5	Ethylbenzene	106	C8H10	000100-41-4	90





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ALS Vial : 49 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0VE2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

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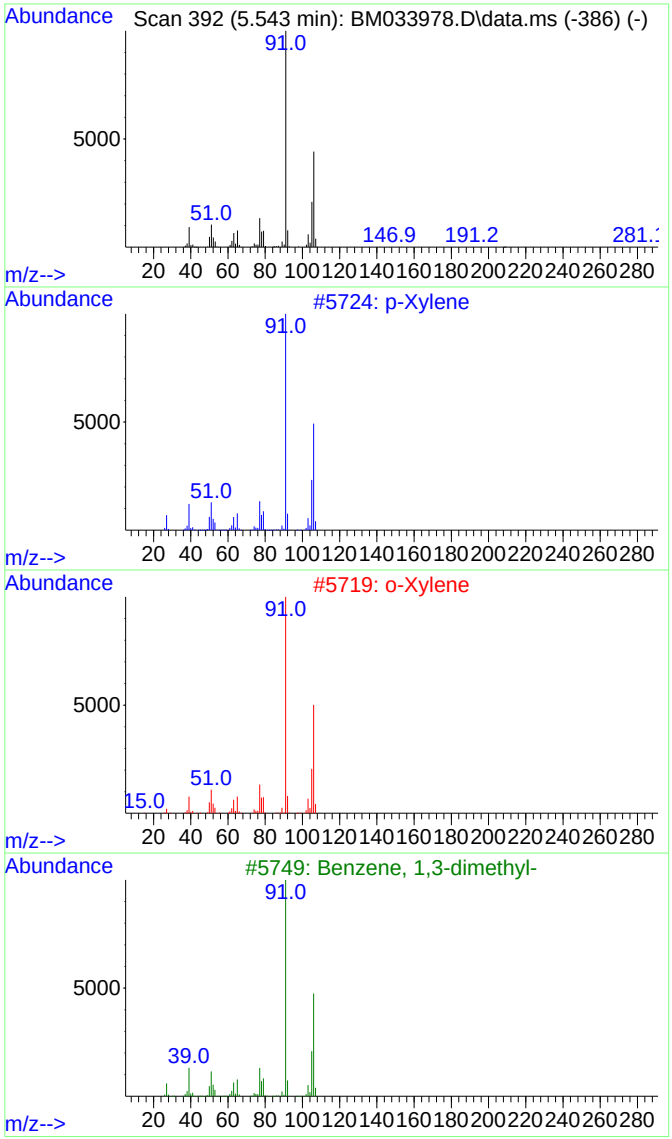
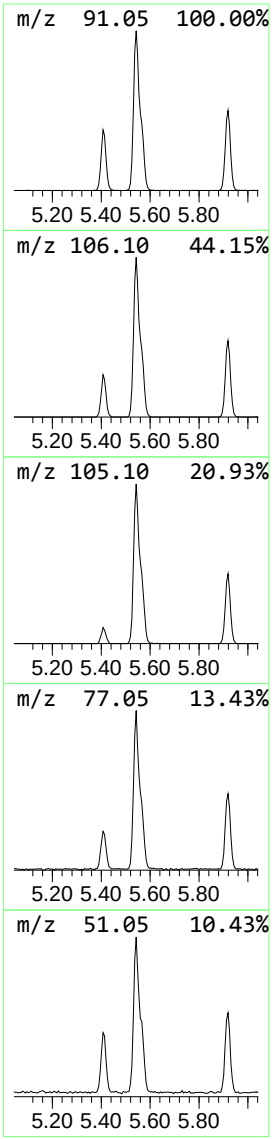
Peak Number 7 p-Xylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.543	36.08 ng/ul	1194880	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	97
2	o-Xylene	106	C8H10	000095-47-6	97
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97





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Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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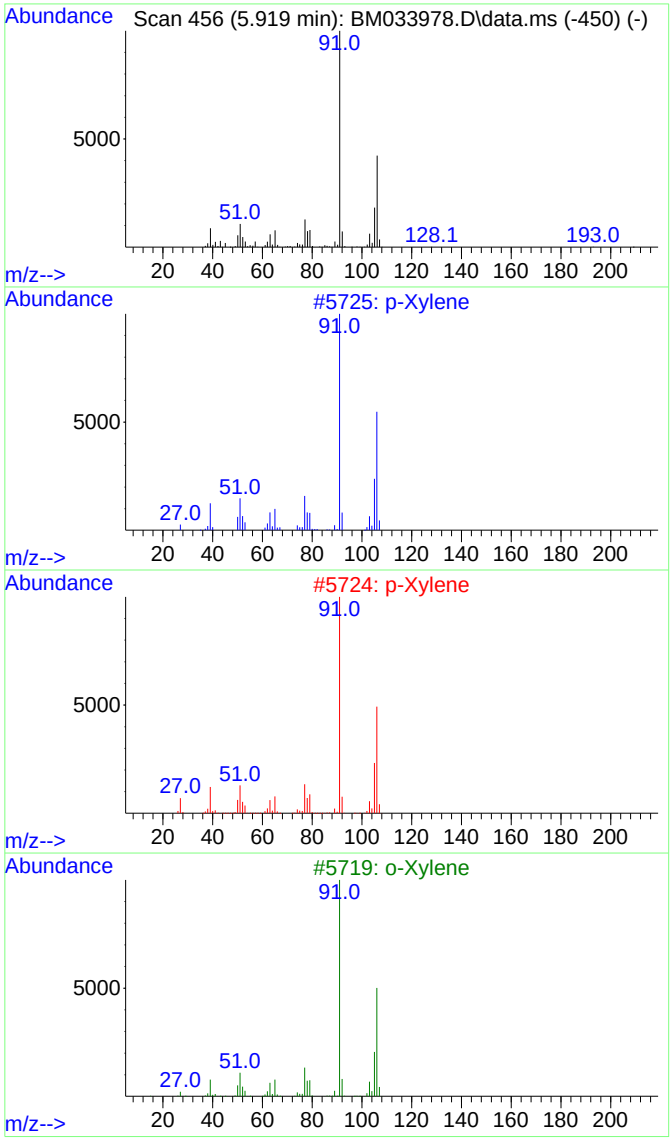
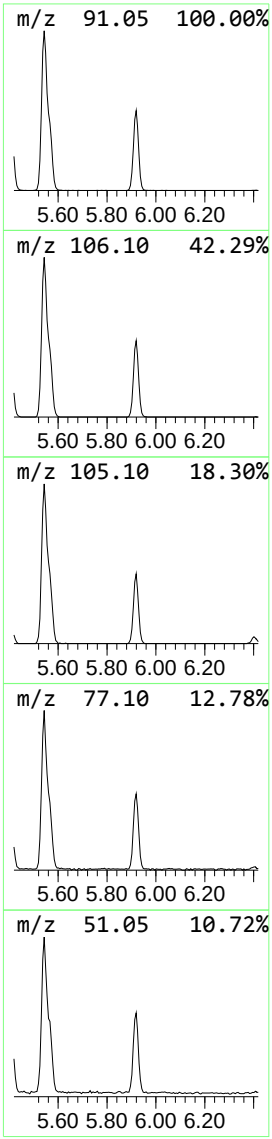
Peak Number 8 o-Xylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.919	14.64 ng/ul	484943	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	o-Xylene	106	C8H10	000095-47-6	97
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97





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Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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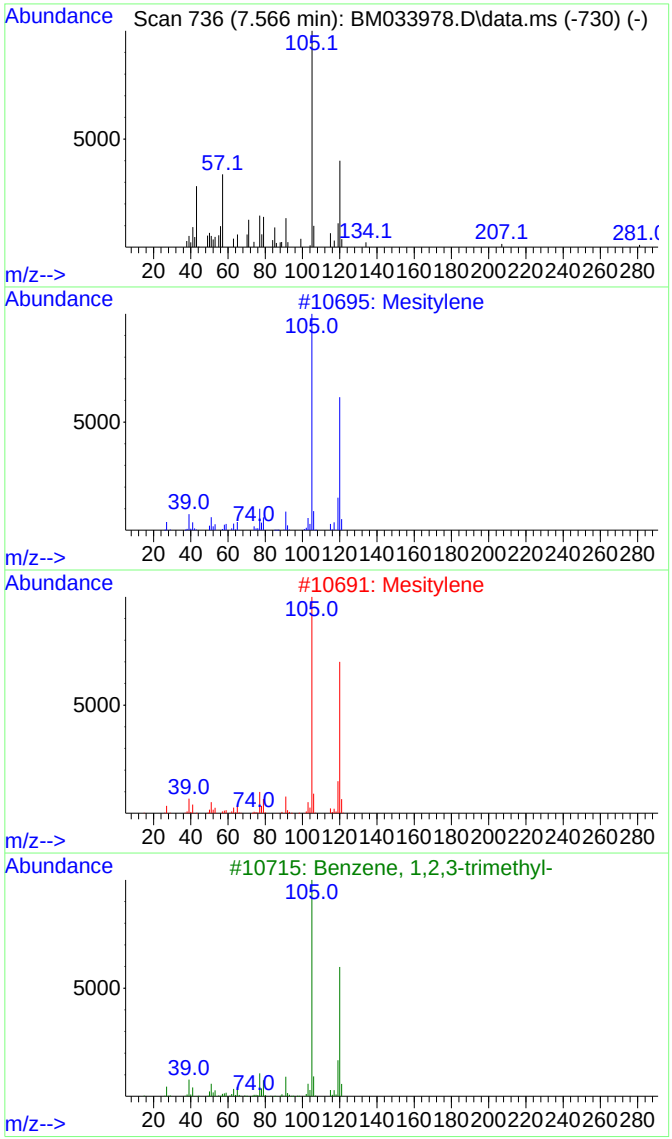
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

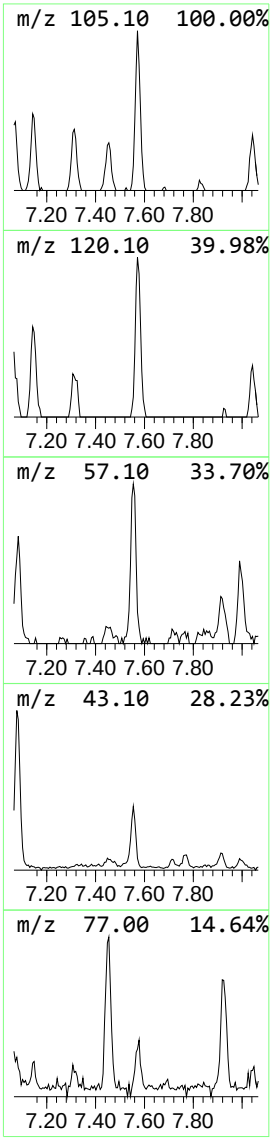
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Mesitylene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.566	2.26 ng/ul	74937	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	81
2	Mesitylene	120	C9H12	000108-67-8	81
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	81
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81





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Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

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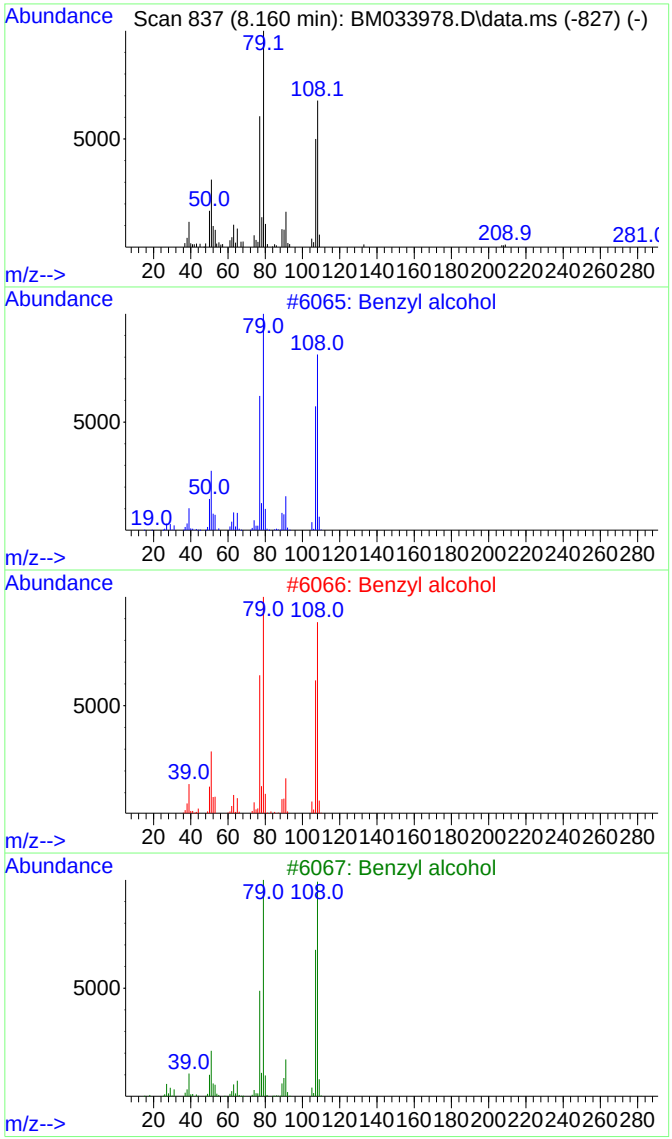
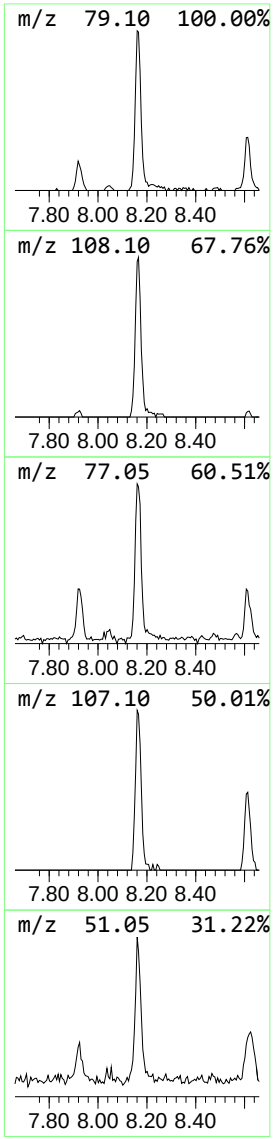
Peak Number 10 Benzyl alcohol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.160	3.47 ng/ul	114941	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	96
2	Benzyl alcohol	108	C7H8O	000100-51-6	96
3	Benzyl alcohol	108	C7H8O	000100-51-6	95
4	Benzyl alcohol	108	C7H8O	000100-51-6	95
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91





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Operator : CG/JU
Sample : N1355-01
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ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

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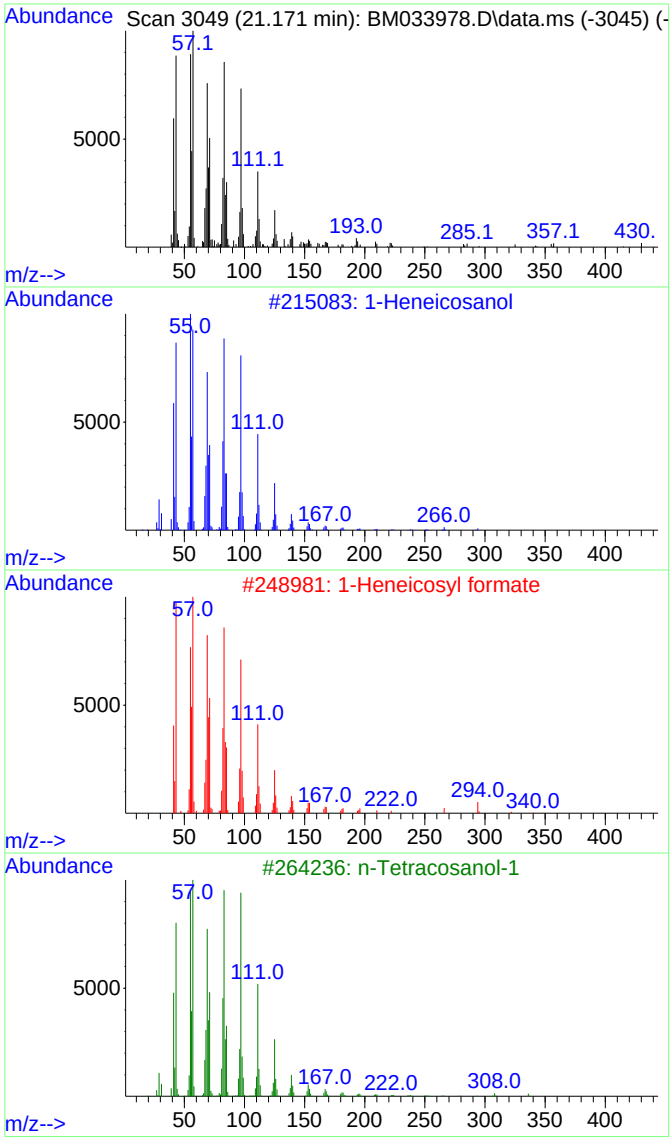
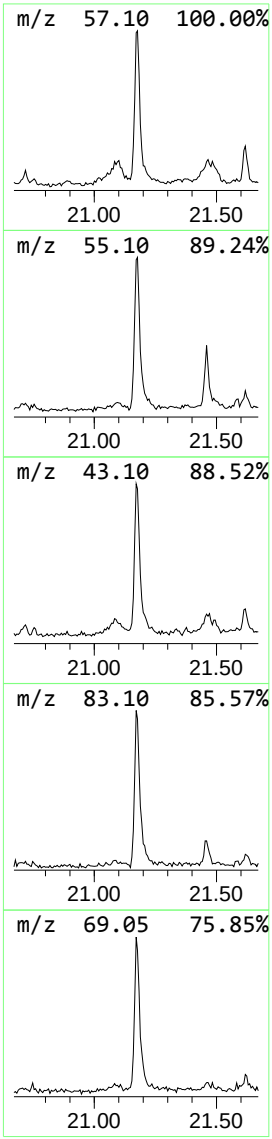
Peak Number 11 1-Heneicosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.171	2.69 ng/ul	245039	Chrysene-d12	21.459

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94





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Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

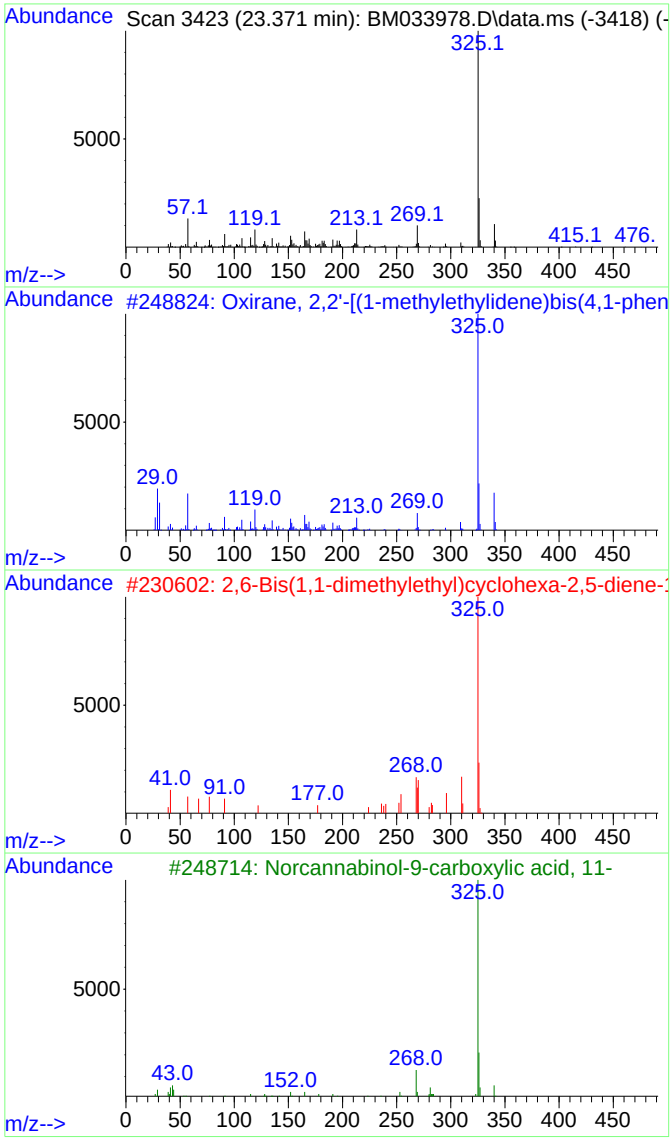
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

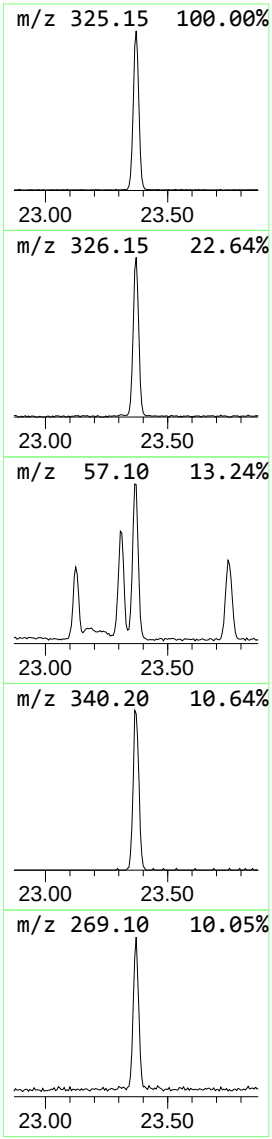
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Oxirane, 2,2'-[(1-methyleth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.371	11.50 ng/ul	823712	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	91
2	2,6-Bis(1,1-dimethylethyl)cycloh...	325	C21H27NO2	017119-01-6	64
3	Norcannabinol-9-carboxylic acid,...	340	C21H24O4	053989-32-5	64
4	3(2H)-Benzofuranone, 2-(hydroxyp...	340	C19H20O4Si	1000503-53-4	64
5	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	64





Tentatively Identified Compound (LSC) summary

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Operator : CG/JU
Sample : N1355-01
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VE2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.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
Methyl Isobutyl...	3.643	37.2	ng/ul	1231480	1	7.925	662286	20.0
2-Pentanol, 4-m...	3.872	3.4	ng/ul	111439	1	7.925	662286	20.0
Butanoic acid	3.954	3.9	ng/ul	128811	1	7.925	662286	20.0
Toluene	4.054	3.6	ng/ul	118005	1	7.925	662286	20.0
2-Pentanone, 4-...	5.001	16.4	ng/ul	543669	1	7.925	662286	20.0
Ethylbenzene	5.407	8.5	ng/ul	280565	1	7.925	662286	20.0
p-Xylene	5.543	36.1	ng/ul	1194880	1	7.925	662286	20.0
o-Xylene	5.919	14.6	ng/ul	484943	1	7.925	662286	20.0
Mesitylene	7.566	2.3	ng/ul	74937	1	7.925	662286	20.0
Benzyl alcohol	8.160	3.5	ng/ul	114941	1	7.925	662286	20.0
1-Heneicosanol	21.171	2.7	ng/ul	245039	5	21.459	1819850	20.0
Oxirane, 2,2'-[...	23.371	11.5	ng/ul	823712	6	23.847	1432130	20.0