LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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: BM033984.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	21	rVB2	36534	58487	0.20%	0.079%
2	3.596	55	61	64	rBV2	29775	45279	0.16%	0.061%
3	3.649	64	70	79	rVV	4636630	5556000	19.10%	7.545%
4	3.719	79	82	96	rVB	140811	208004	0.71%	0.282%
5	3.872	104	108	115	rVB	62584	79881	0.27%	0.108%
6	3.990	115	128	132	rVV	123464	330995	1.14%	0.450%
7	4.054	134	139	149	rVB	106568	149103	0.51%	0.202%
8	5.001	294	300	313	rBV	246192	340175	1.17%	0.462%
9	5.407	362	369	378	rVB	609006	865811	2.98%	1.176%
10	5.543	385	392	403	rBV	2586249	4392463	15.10%	5.965%
11	5.748	423	427	433	rBV	35923	47320	0.16%	0.064%
12	5.919	450	456	463	rBV	913969	1348155	4.63%	1.831%
13	5.984	463	467	474	rVB	278424	397012	1.36%	0.539%
14	6.913	617	625	631	rBV5	13116	31248	0.11%	0.042%
15	7.078	646	653	662	rVV	561888	932553	3.21%	1.266%
16	7.248	675	682	689	rVV	420431	652211	2.24%	0.886%
17	7.448	709	716	725	rBV	559920	893375	3.07%	1.213%
18	7.572	730	737	743	rVB2	30385	71093	0.24%	0.097%
19	7.919	790	796	804	rBV	465720	754178	2.59%	1.024%
20	8.483	885	892	896	rBV2	21664	36520	0.13%	0.050%
21	8.631	909	917	926	rBV	682640	1092049	3.75%	1.483%
22	8.748	932	937	943	rVB	21738	36431	0.13%	0.049%
23	9.083	988	994	1002	rVB	532974	841916	2.89%	1.143%
24	9.530	1065	1070	1078	rBV	27856	45086	0.15%	0.061%
25	9.813	1104	1118	1124	rBV	438217	732698	2.52%	0.995%
26	10.354	1202	1210	1222	rBV	686036	1171776	4.03%	1.591%
27	10.513	1232	1237	1245	rBV	58388	92926	0.32%	0.126%
28	10.736	1266	1275	1281	rBV	631220	1089149	3.74%	1.479%
29	10.866	1290	1297	1311	rBV	500674	879138	3.02%	1.194%
30	13.395	1720	1727	1734	rBV4	36824	72673	0.25%	0.099%
31	13.971	1819	1825	1830	rBV	1001446	1432231	4.92%	1.945%
32	14.071	1835	1842	1844	rVV3	18858	40744	0.14%	0.055%
33	14.107	1844	1848	1856	rVB3	19939	35180	0.12%	0.048%
34	14.254	1867	1873	1886	rVB	1176719	1744611	6.00%	2.369%
35	14.560	1914	1925	1936	rVB	907116	1356171	4.66%	1.842%
36	14.748	1948	1957	1972	rBV	297800	563968	1.94%	0.766%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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	15.548	2087	2093	2103	rBV	1470917	2056877	7.07%	2.793%
38	15.659	2106	2112	2120	rBV	419882	570051	1.96%	0.774%
39	17.295	2384	2390	2399	rBV	986650	1399937	4.81%	1.901%
40	17.395	2400	2407	2419	rVB	1449169	2045416	7.03%	2.778%
41	18.189	2538	2542	2550	rBV	66861	111015	0.38%	0.151%
42	18.506	2591	2596	2602	rBV	73435	100216	0.34%	0.136%
43	19.465	2754	2759	2764	rBV6	26395	45520	0.16%	0.062%
44	19.677	2789	2795	2800	rBV	1622795	2219511	7.63%	3.014%
45	19.718	2800	2802	2809	rVB	93874	134426	0.46%	0.183%
46	21.077	3030	3033	3037	rBV2	49304	68607	0.24%	0.093%
47	21.171	3045	3049	3058	rVB2	188942	277163	0.95%	0.376%
48	21.459	3093	3098	3103	rBV	1145447	1479396	5.08%	2.009%
49	21.506	3103	3106	3112	rVB	212889	272014	0.93%	0.369%
50	22.153	3212	3216	3222	rVB2	359661	469042	1.61%	0.637%
51	23.306	3408	3412	3416	rBV	160451	248242	0.85%	0.337%
52	23.394	3416	3427	3436	rVB	14831705	29094979	100.00%	39.513%
53	23.694	3471	3478	3484	rBV2	1116844	2187092	7.52%	2.970%
54	23.847	3498	3504	3516	rVB2	761251	1517843	5.22%	2.061%
55	24.630	3632	3637	3644	rVB	223386	478367	1.64%	0.650%
56	24.847	3670	3674	3684	rVB	215253	442284	1.52%	0.601%

Sum of corrected areas: 73634608

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Data File : BM033984.D
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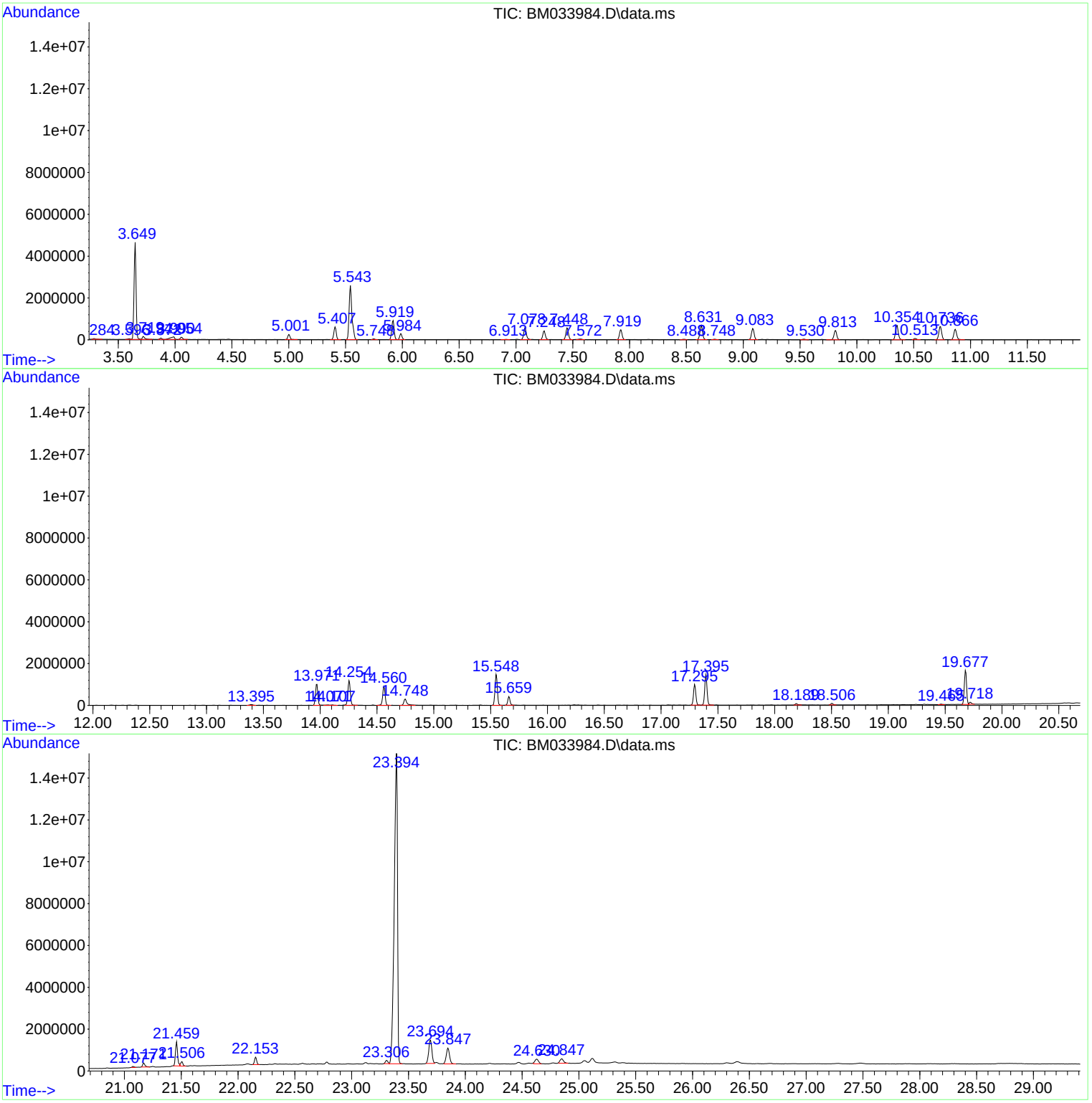
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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 : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
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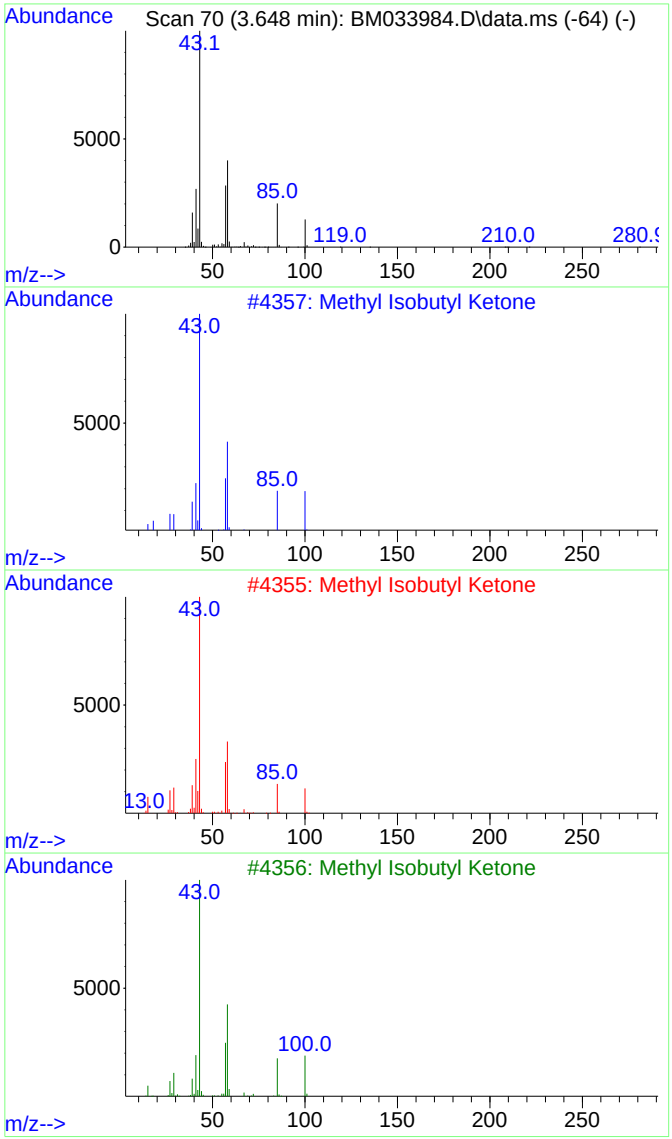
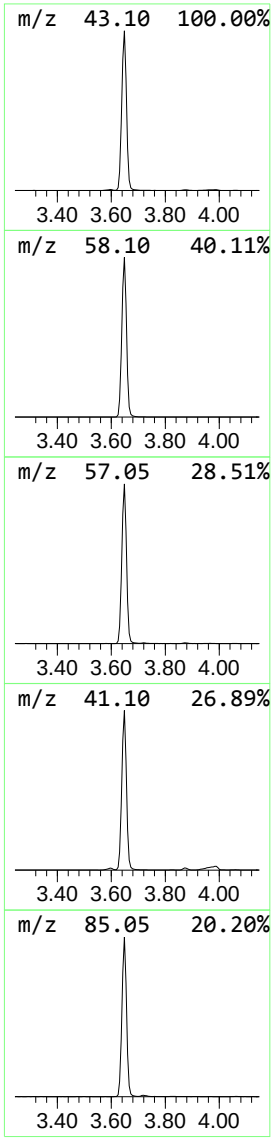
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.648	147.34 ng/ul	5556000	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	86
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	83
5	2-Hexanone	100	C6H12O	000591-78-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

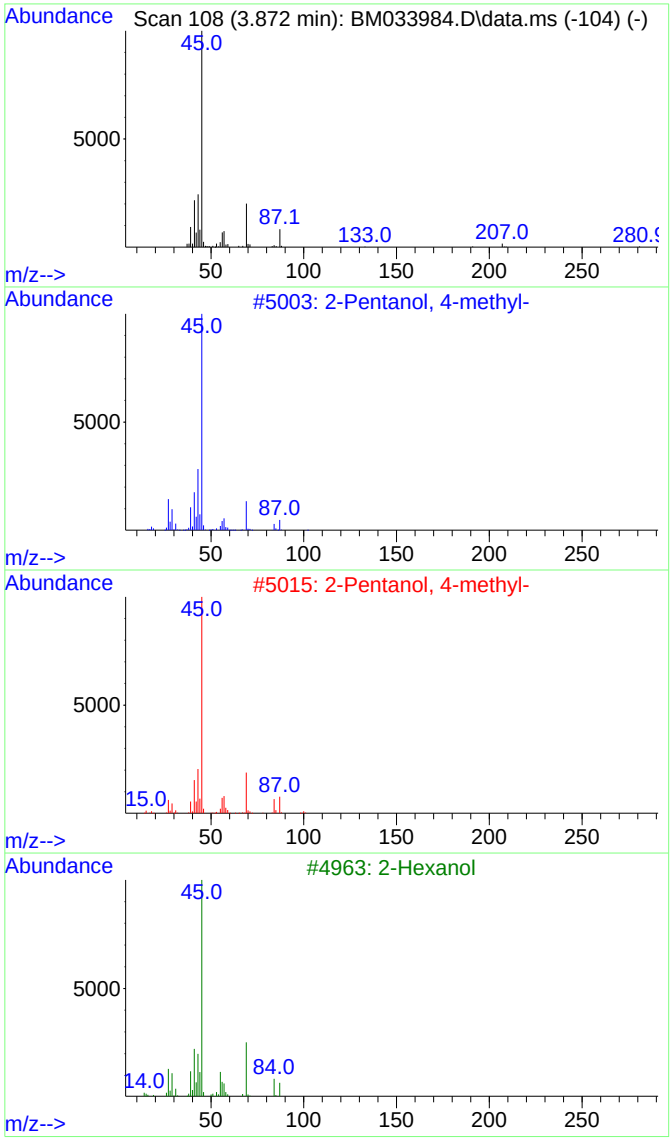
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

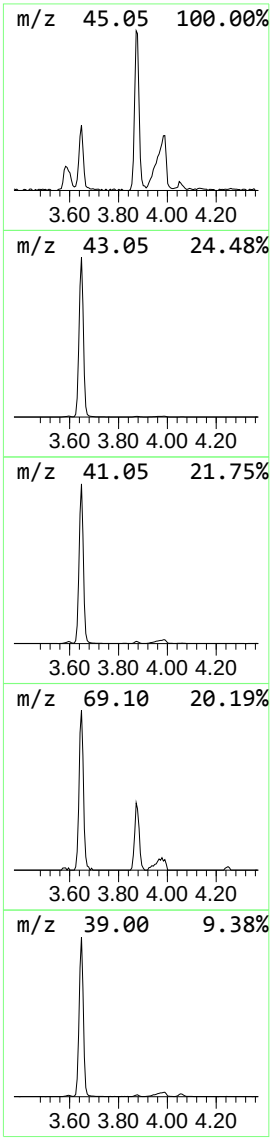
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanol, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.872	2.12 ng/ul	79881	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	78
3	2-Hexanol	102	C6H14O	000626-93-7	72
4	2-Hexanol	102	C6H14O	000626-93-7	72
5	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

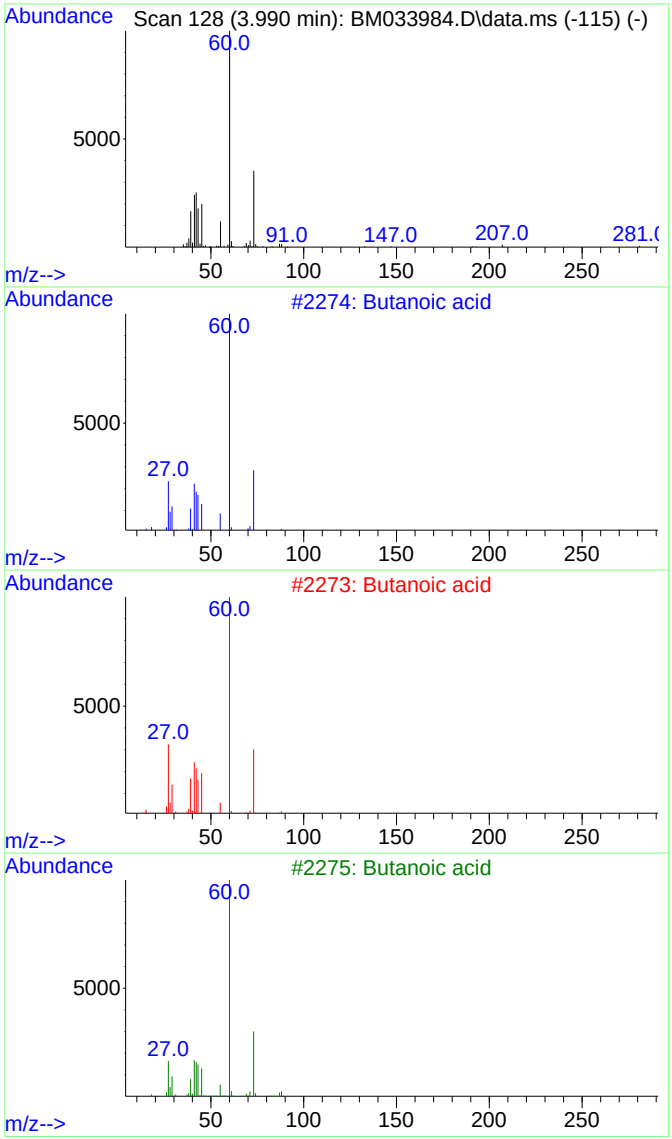
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

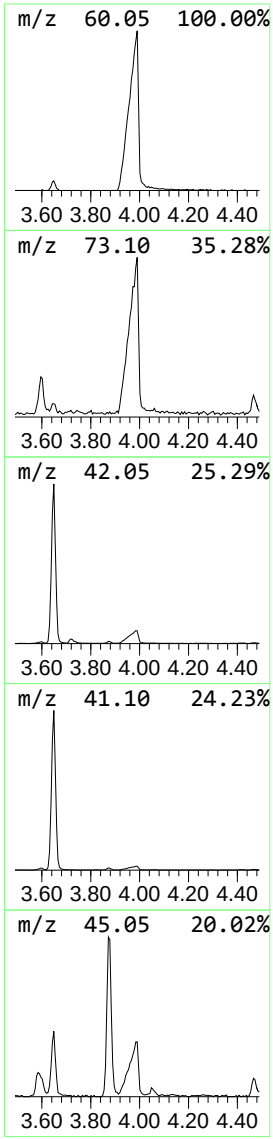
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.990	8.78 ng/ul	330995	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	83
2	Butanoic acid	88	C4H8O2	000107-92-6	80
3	Butanoic acid	88	C4H8O2	000107-92-6	72
4	Butanoic acid	88	C4H8O2	000107-92-6	64
5	Hexanoic acid	116	C6H12O2	000142-62-1	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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

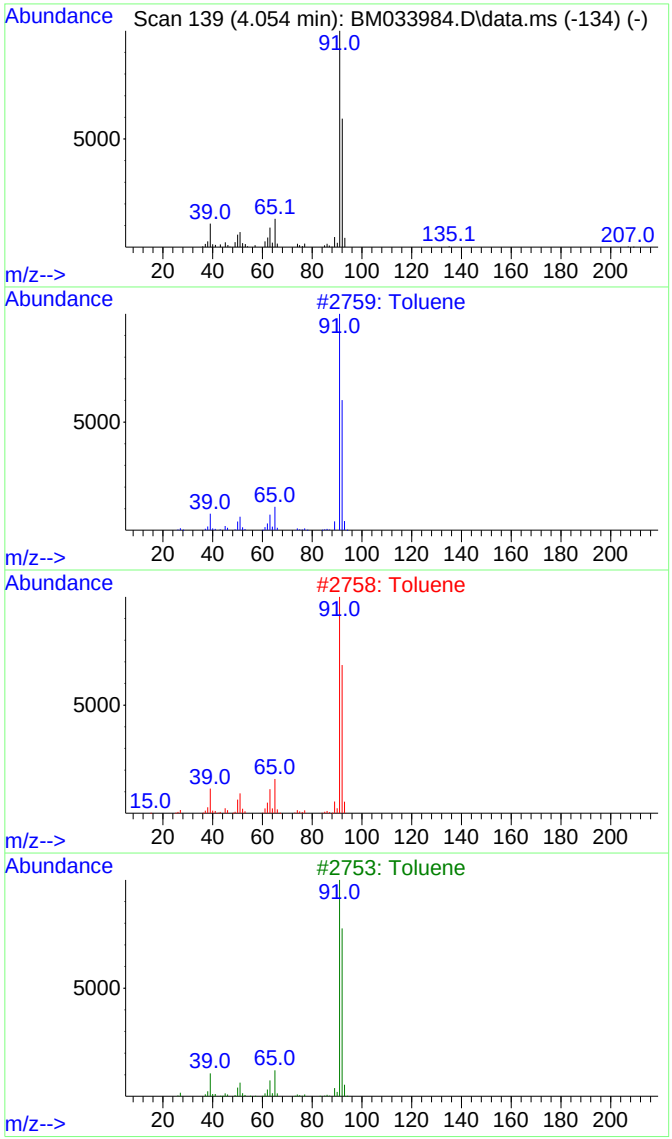
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

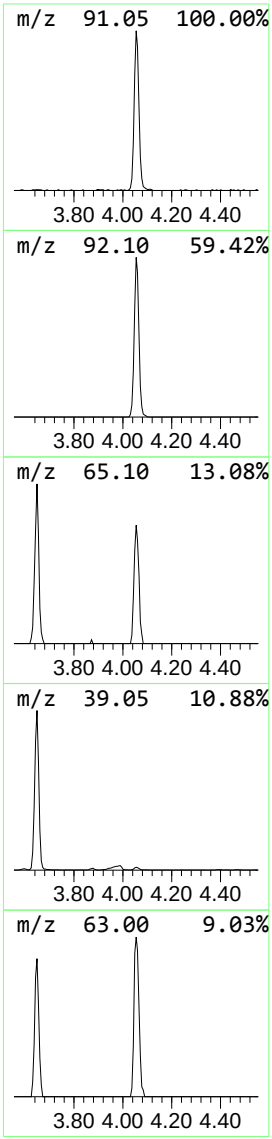
****

Peak Number 4 Toluene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.054	3.95 ng/ul	149103	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	93
2	Toluene	92	C7H8	000108-88-3	93
3	Toluene	92	C7H8	000108-88-3	91
4	Toluene	92	C7H8	000108-88-3	90
5	Toluene	92	C7H8	000108-88-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

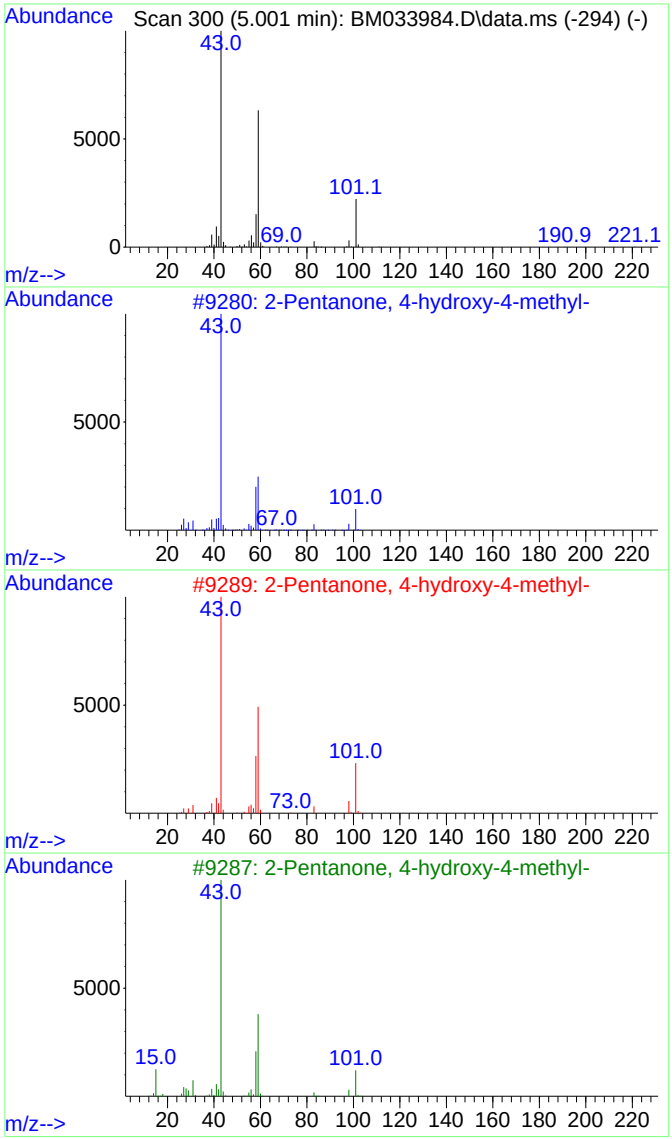
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

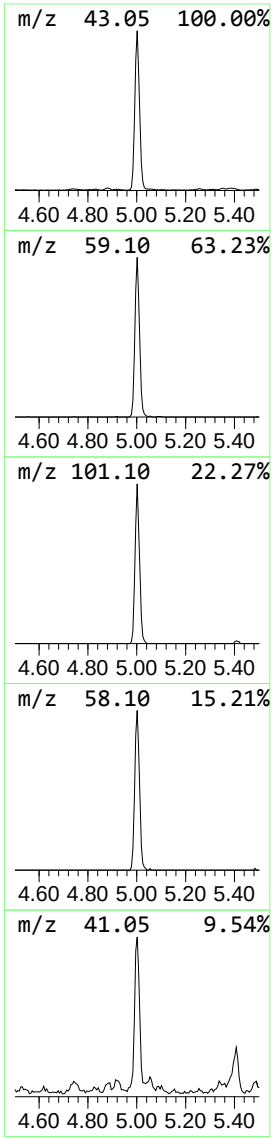
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.001	9.02 ng/ul	340175	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	72
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	56
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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

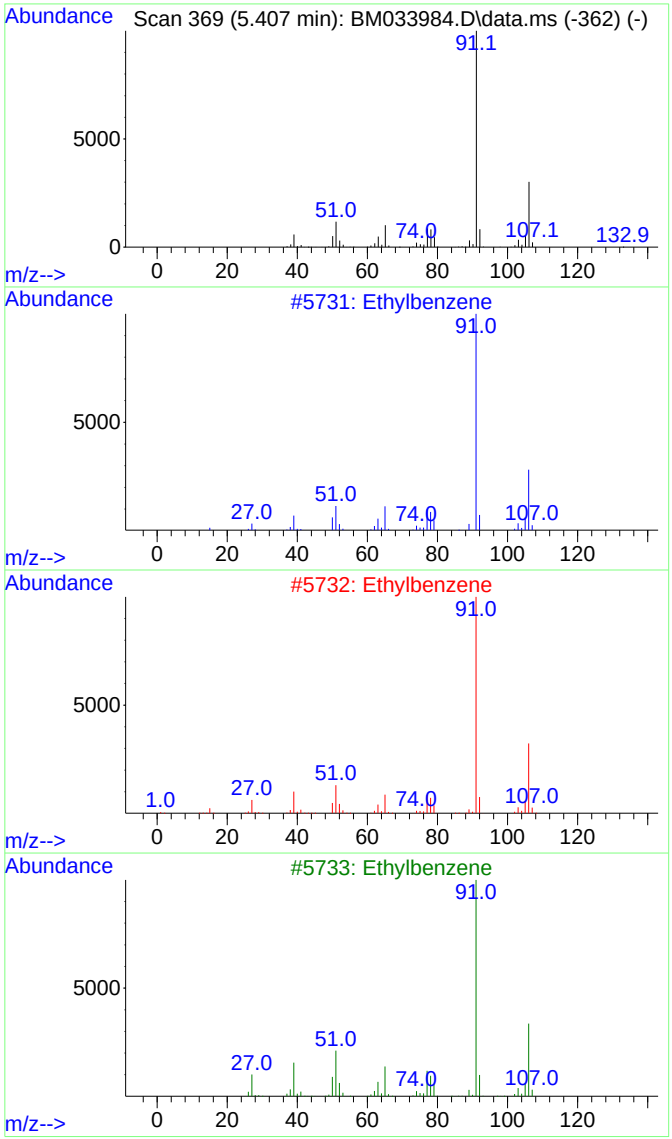
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

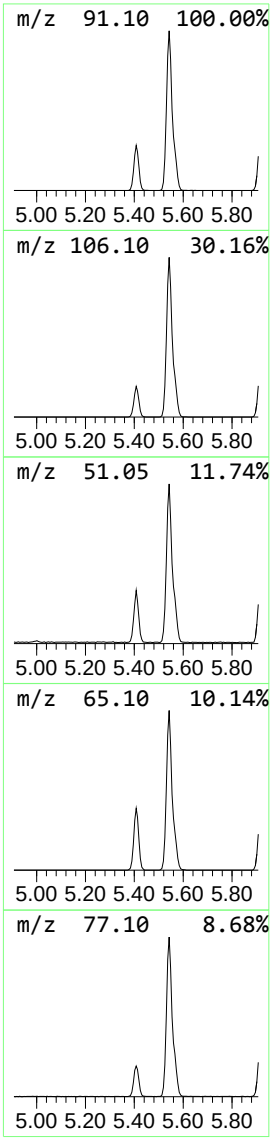
****

Peak Number 6 Ethylbenzene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.407	22.96 ng/ul	865811	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	91
4	Ethylbenzene	106	C8H10	000100-41-4	90
5	Ethylbenzene	106	C8H10	000100-41-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

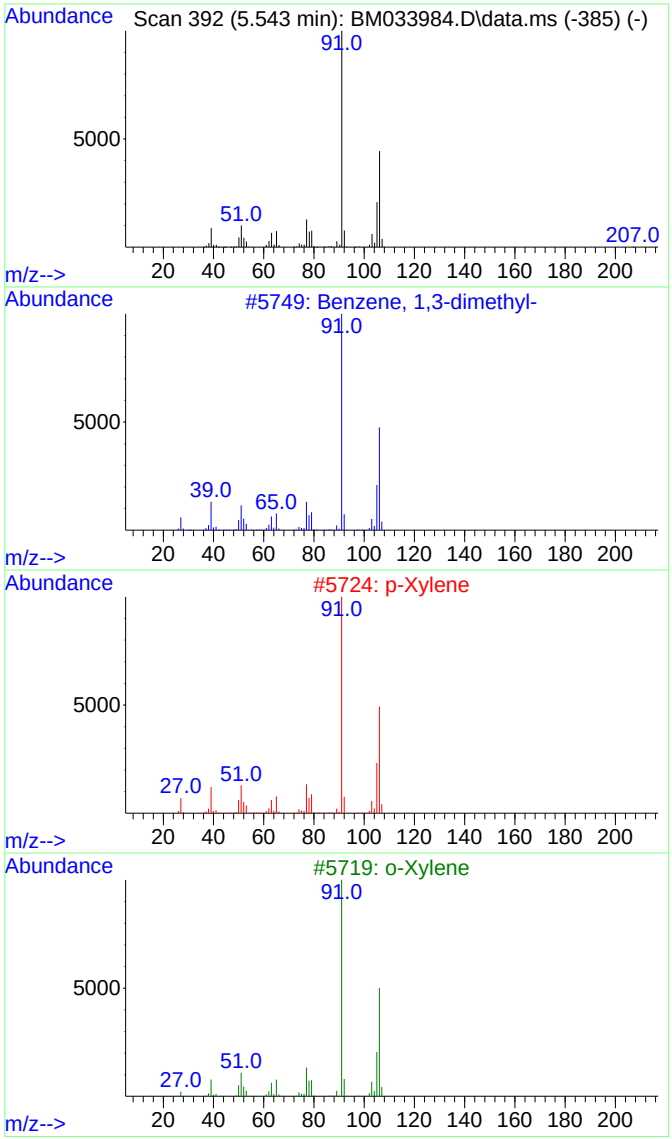
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

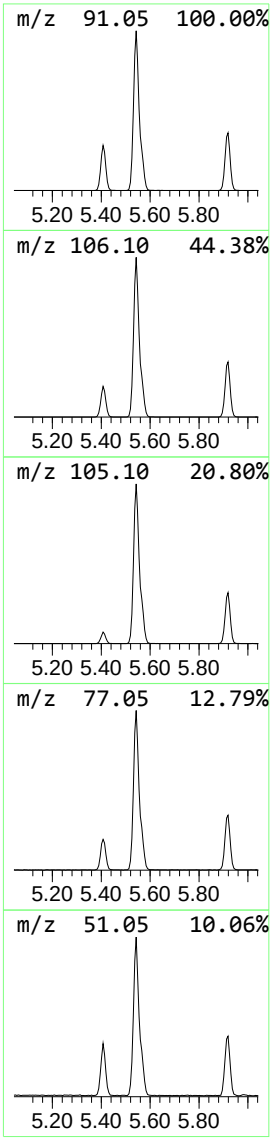
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.543	116.48 ng/ul	4392460	1,4-Dichlorobenzene-d4	7.925

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0VF4

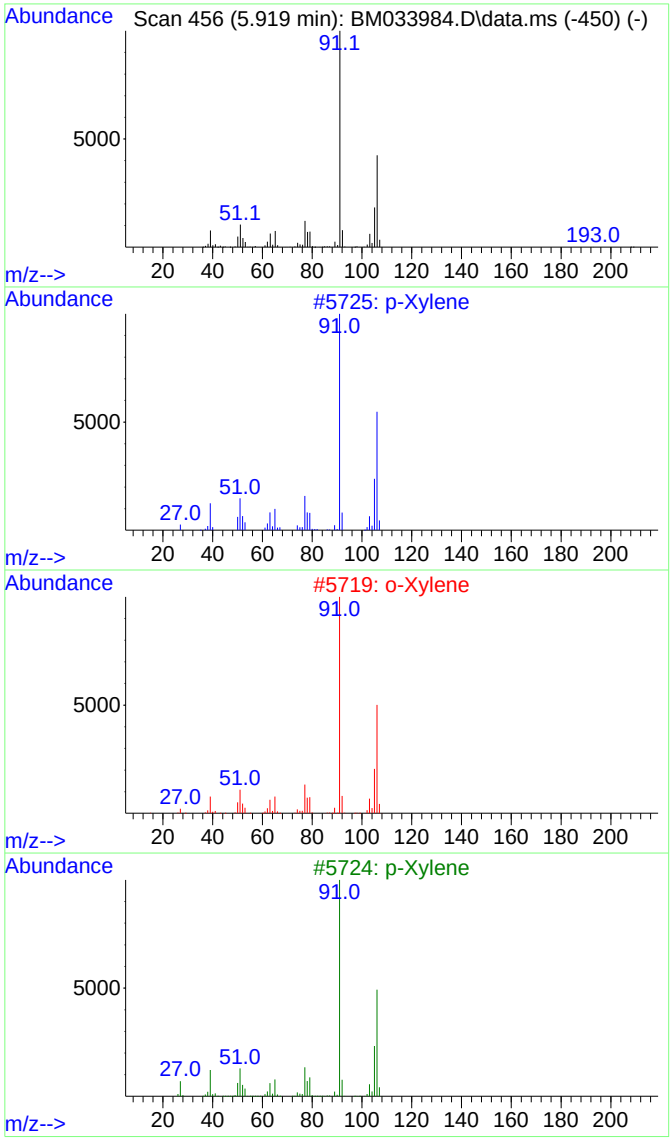
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

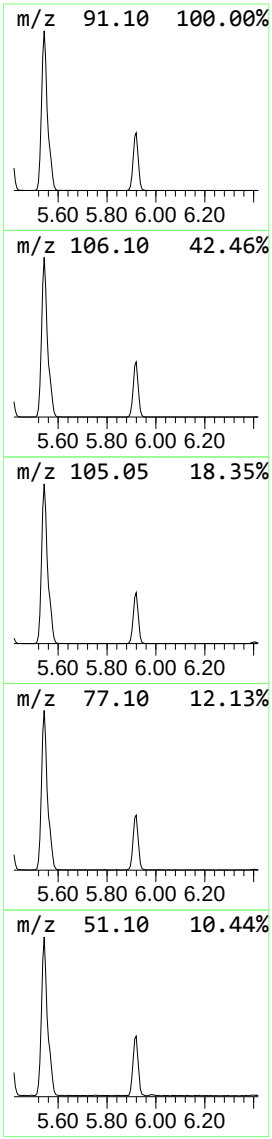
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 p-Xylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.919	35.75 ng/ul	1348160	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	p-Xylene	106	C8H10	000106-42-3	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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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

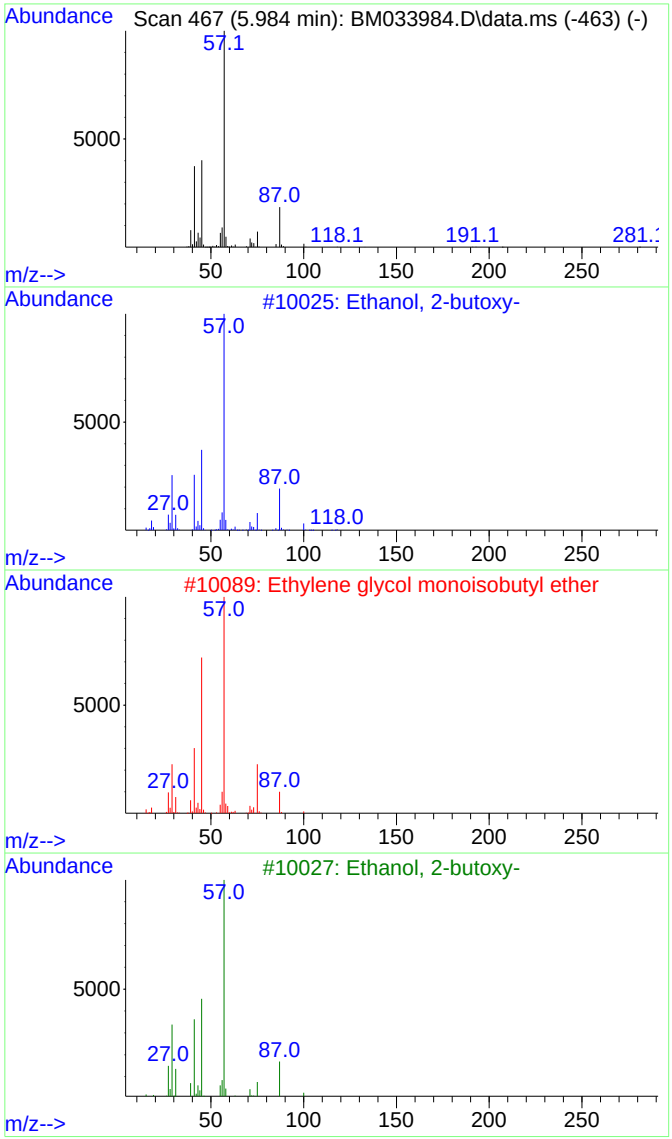
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

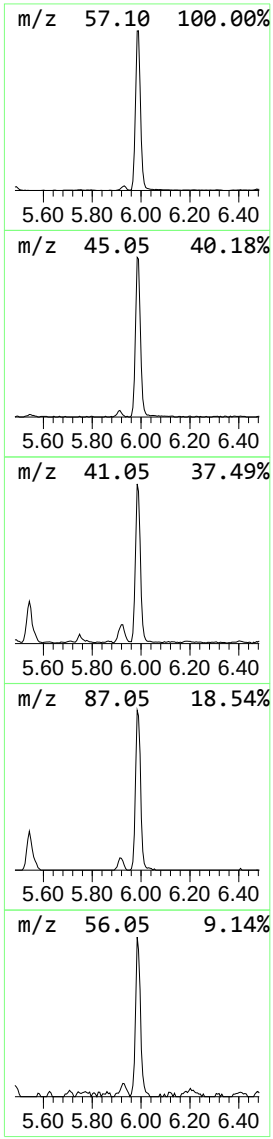
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Ethanol, 2-butoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.984	10.53 ng/ul	397012	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
2	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
4	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	59





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

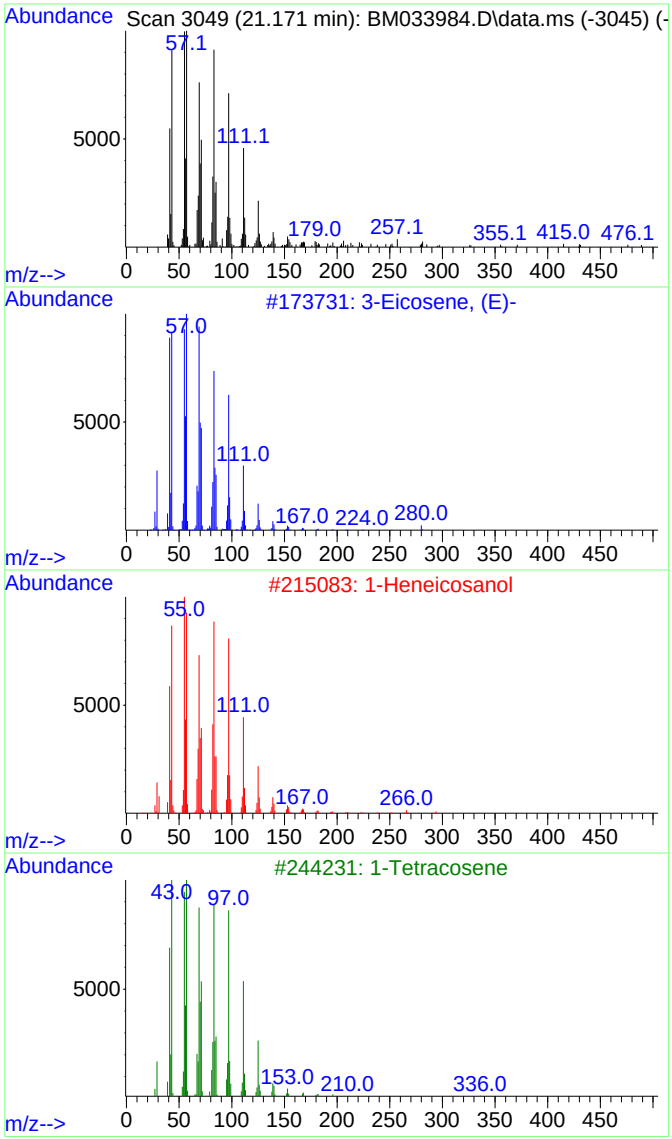
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

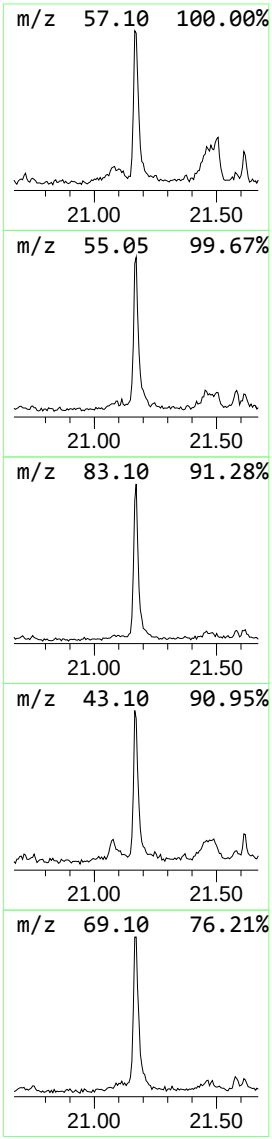
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.171	3.75 ng/ul	277163	Chrysene-d12	21.459

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	1-Tetracosene	336	C24H48	010192-32-2	95
4	Nonacos-1-ene	406	C29H58	018835-35-3	94
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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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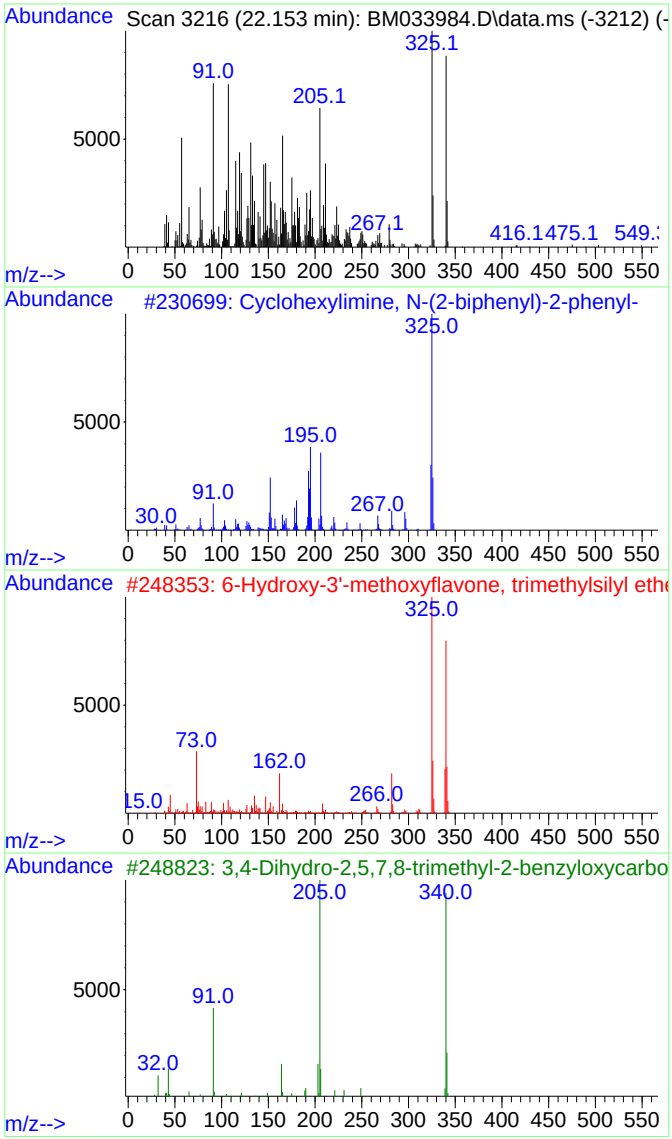
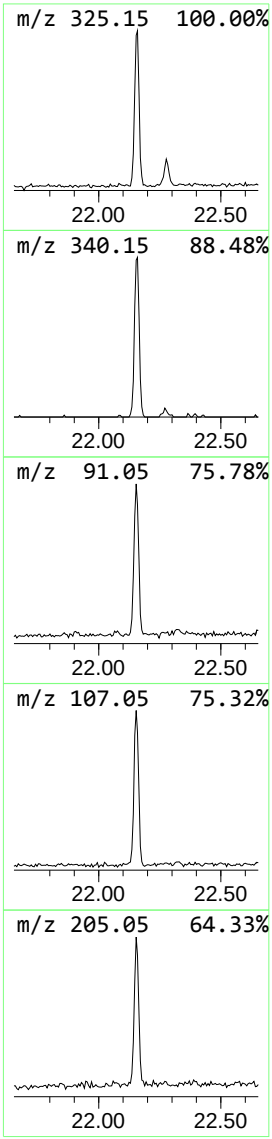
Peak Number 11 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.153	6.34 ng/ul	469042	Chrysene-d12	21.459

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexylimine, N-(2-biphenyl)-...	325	C24H23N	1000161-05-7	45
2	6-Hydroxy-3'-methoxyflavone, tri...	340	C19H20O4Si	1000448-97-9	42
3	3,4-Dihydro-2,5,7,8-trimethyl-2-...	340	C21H24O4	1000432-38-4	35
4	3-(3',4'-Dimethoxyphenyl)-6-etho...	340	C20H20O5	720674-30-6	25
5	3-(4-Hydroxy-3,5-dimethoxybenzyl...	325	C19H19NO4	1000504-86-5	18





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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

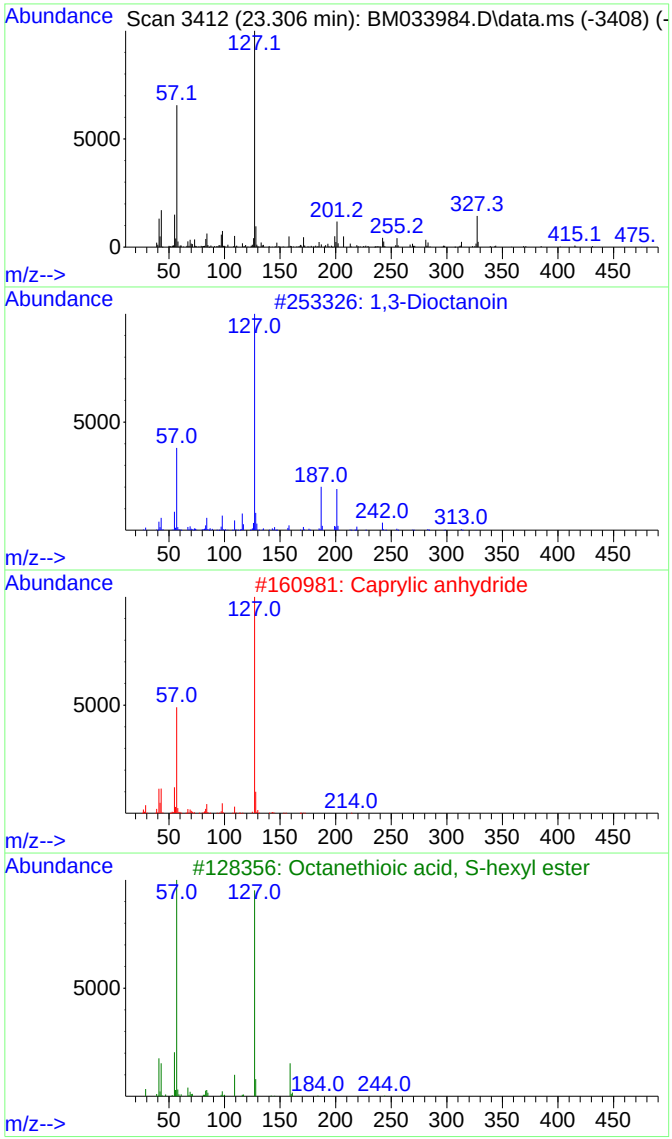
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

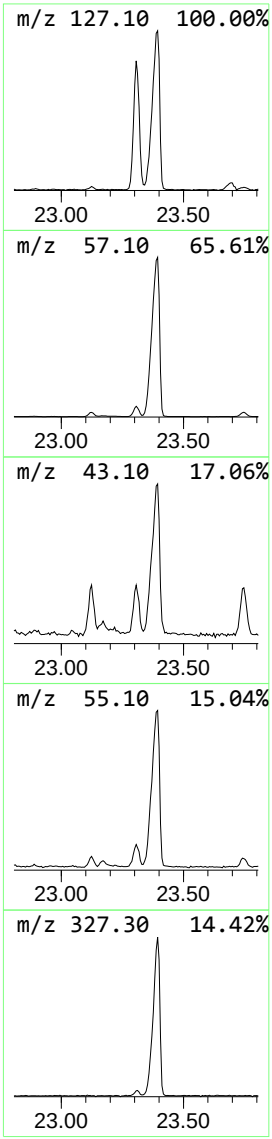
****

Peak Number 12 1,3-Dioctanoin Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.306	3.27 ng/ul	248242	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioctanoin	344	C19H36O5	001429-66-9	59
2	Caprylic anhydride	270	C16H30O3	000623-66-5	58
3	Octanethioic acid, S-hexyl ester	244	C14H28OS	055590-85-7	58
4	L-Valine, N-dimethylaminomethyle...	200	C10H20N2O2	1000375-63-8	53
5	3-Fluoro-5-methoxypyridine	127	C6H6FNO	1060801-62-8	53





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

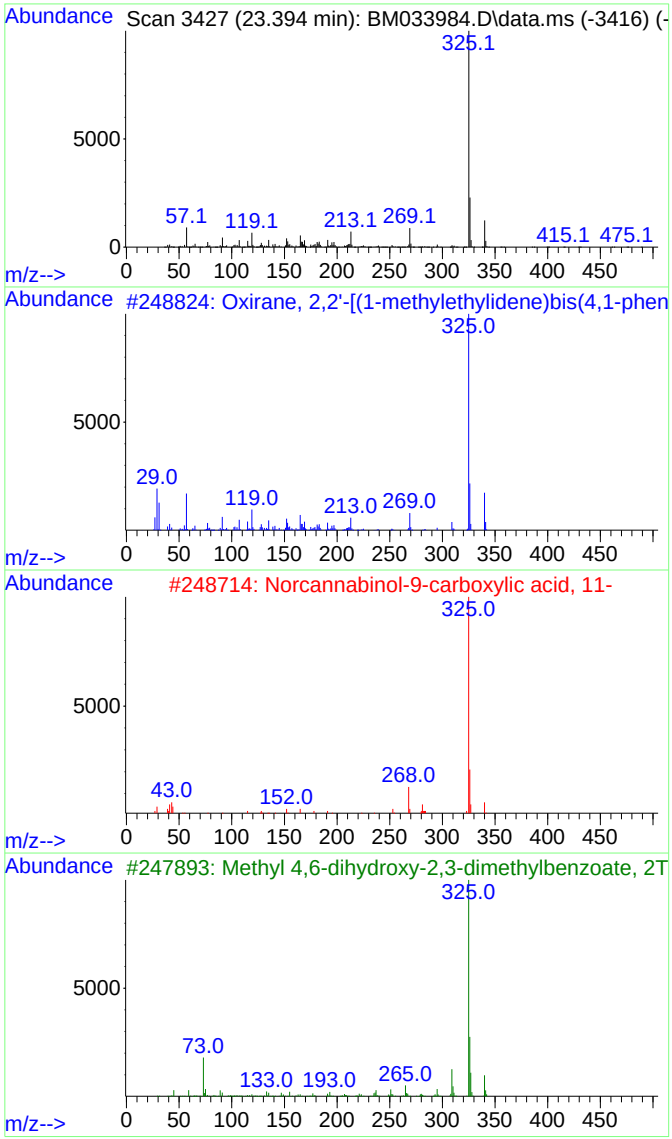
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

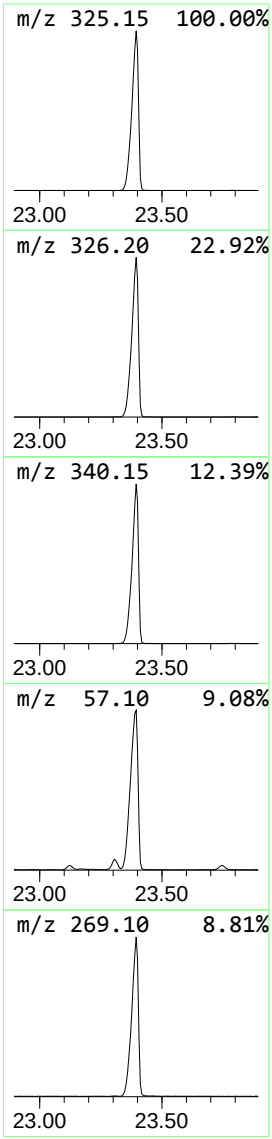
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Oxirane, 2,2'-[(1-methyleth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.394	383.37 ng/ul	29095000	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	89
2	Norcannabinol-9-carboxylic acid,...	340	C21H24O4	053989-32-5	80
3	Methyl 4,6-dihydroxy-2,3-dimethy...	340	C16H28O4Si2	1000502-58-9	72
4	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	70
5	3(2H)-Benzofuranone, 2-(hydroxyp...	340	C19H20O4Si	1000503-53-4	56





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

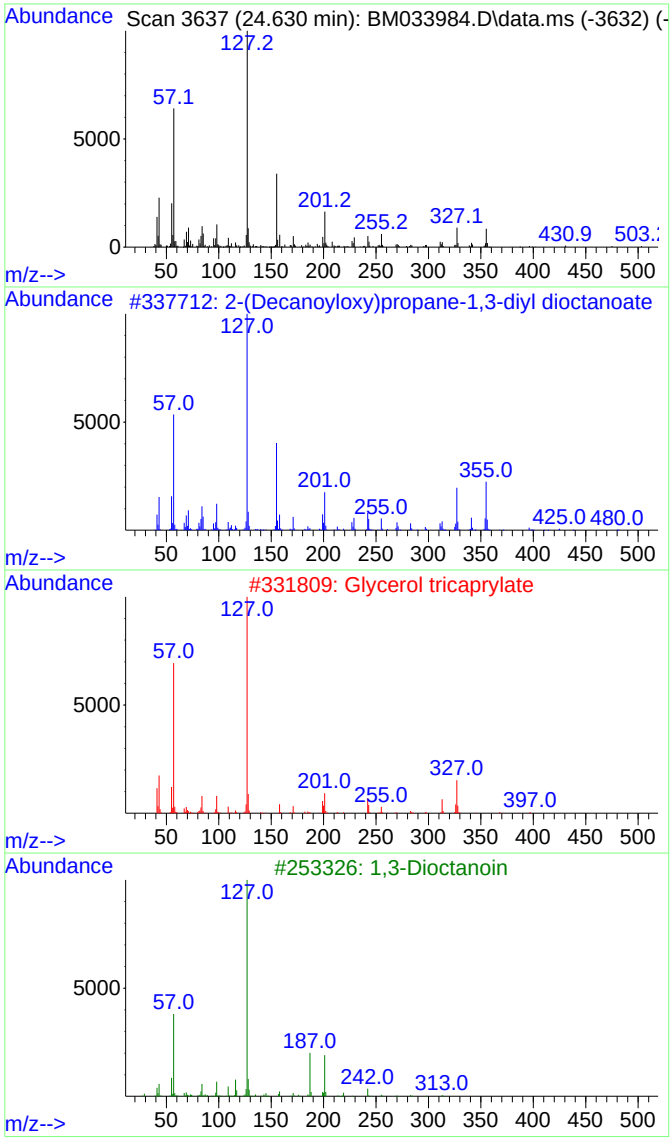
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

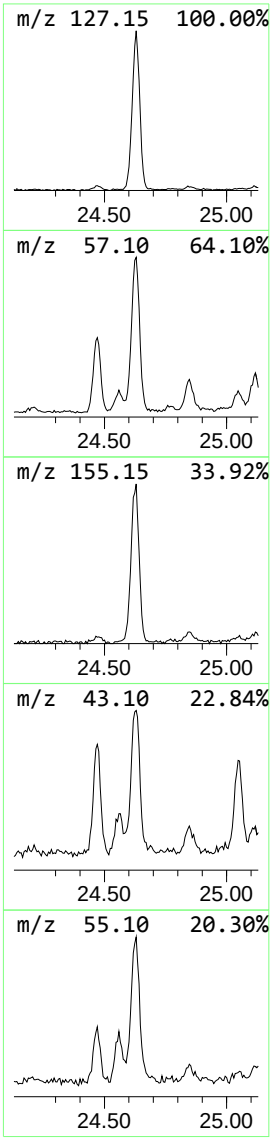
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 2-(Decanoyloxy)propane-1,3-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.630	6.30 ng/ul	478367	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(Decanoyloxy)propane-1,3-diyl ...	498	C29H54O6	033368-87-5	91
2	Glycerol tricaprylate	470	C27H50O6	000538-23-8	59
3	1,3-Dioctanoin	344	C19H36O5	001429-66-9	53
4	1,2-Dioctanoyl-sn-glycerol, O-ac...	386	C21H38O6	1000452-09-1	53
5	5-Bromo-4'-chlorosalicylanilide	325	C13H9BrClNO2	003679-64-9	52





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Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF4

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

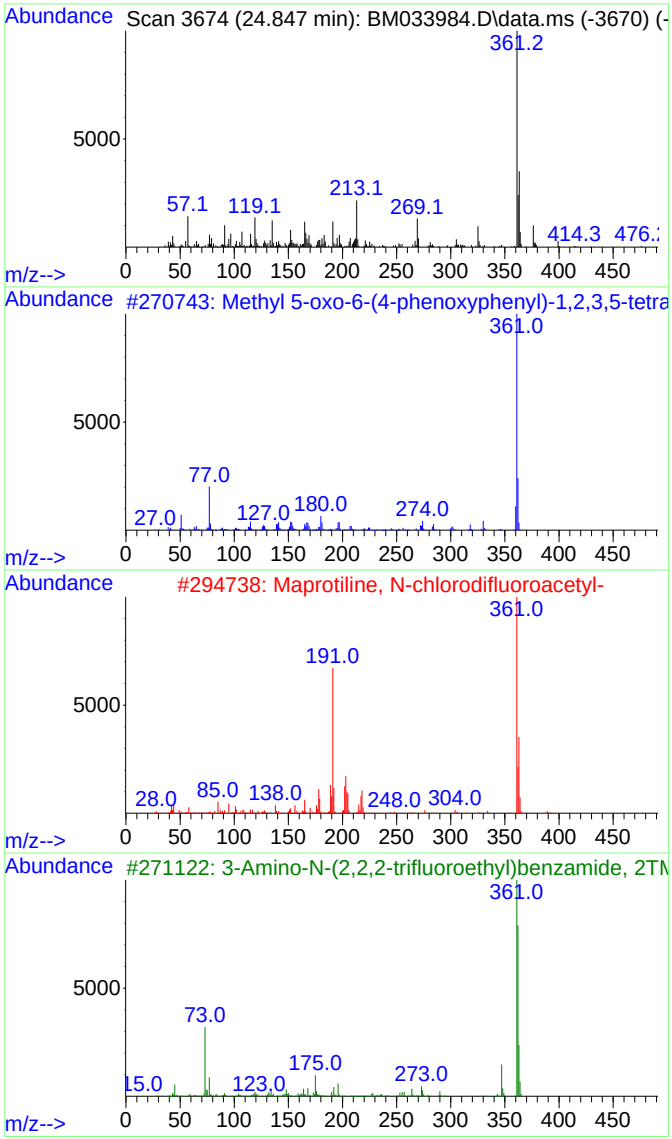
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

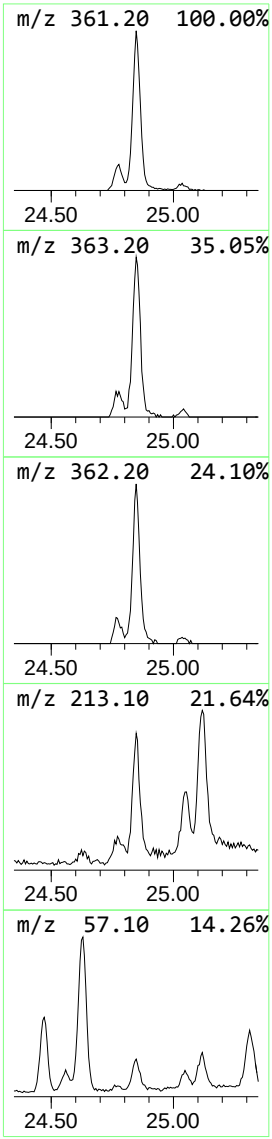
****

Peak Number 15 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.847	5.83 ng/ul	442284	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 5-oxo-6-(4-phenoxyphenyl)...	361	C22H19NO4	1000482-02-4	49
2	Maprotiline, N-chlorodifluoroace...	389	C22H22ClF2NO	1000448-25-9	42
3	3-Amino-N-(2,2,2-trifluoroethyl)...	362	C15H25F3N2OSi2	1000501-27-9	38
4	Thiazolo[3,2-b]1,2,4-triazole, 2...	361	C16H10F3N5S	266679-25-8	38
5	Benzo[b,f]-1,2,4-triazolo[4,3-d]...	361	C22H11N5O	1000277-12-1	38





Tentatively Identified Compound (LSC) summary

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Data File : BM033984.D
Acq On : 02 Feb 2022 21:24
Operator : CG/JU
Sample : N1355-10
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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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Methyl Isobutyl...	3.648	147.3	ng/ul	5556000	1	7.925	754178	20.0
2-Pentanol, 4-m...	3.872	2.1	ng/ul	79881	1	7.925	754178	20.0
Butanoic acid	3.990	8.8	ng/ul	330995	1	7.925	754178	20.0
Toluene	4.054	4.0	ng/ul	149103	1	7.925	754178	20.0
2-Pentanone, 4-...	5.001	9.0	ng/ul	340175	1	7.925	754178	20.0
Ethylbenzene	5.407	23.0	ng/ul	865811	1	7.925	754178	20.0
Benzene, 1,3-di...	5.543	116.5	ng/ul	4392460	1	7.925	754178	20.0
p-Xylene	5.919	35.8	ng/ul	1348160	1	7.925	754178	20.0
Ethanol, 2-butoxy-	5.984	10.5	ng/ul	397012	1	7.925	754178	20.0
(DEL) Alkane: S...	21.171	3.8	ng/ul	277163	5	21.459	1479400	20.0
unknown-01	22.153	6.3	ng/ul	469042	5	21.459	1479400	20.0
1,3-Dioctanoin	23.306	3.3	ng/ul	248242	6	23.847	1517840	20.0
Oxirane, 2,2'-[...	23.394	383.4	ng/ul	29095000	6	23.847	1517840	20.0
2-(Decanoyloxy)...	24.630	6.3	ng/ul	478367	6	23.847	1517840	20.0
unknown-02	24.847	5.8	ng/ul	442284	6	23.847	1517840	20.0