LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

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-BM121523.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM043739.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.145	667	673	681	rVB	76237	136253	0.12%	0.072%
2	7.492	726	732	740	rVB2	84849	143427	0.12%	0.075%
3	7.592	740	749	757	rBV	86817	156302	0.13%	0.082%
4	7.963	805	812	818	rBV	1024015	1665607	1.44%	0.875%
5	8.028	818	823	836	rVB	165611	324599	0.28%	0.170%
6	8.692	930	936	945	rBV2	83773	158284	0.14%	0.083%
7	9.139	1004	1012	1022	rBV	70204	147303	0.13%	0.077%
8	9.863	1127	1135	1142	rBV	57797	117550	0.10%	0.062%
9	10.410	1222	1228	1237	rBV2	71550	160765	0.14%	0.084%
10	10.769	1283	1289	1302	rVB	1246595	2209315	1.91%	1.160%
11	11.998	1494	1498	1508	rVB3	64856	129999	0.11%	0.068%
12	12.192	1513	1531	1549	rBV2	394527	1865441	1.61%	0.979%
13	12.827	1634	1639	1646	rVB	169743	282798	0.24%	0.148%
14	13.210	1695	1704	1707	rBV4	116731	270258	0.23%	0.142%
15	13.386	1728	1734	1742	rBV4	83647	251748	0.22%	0.132%
16	13.598	1766	1770	1774	rBV3	85366	122338	0.11%	0.064%
17	13.857	1808	1814	1819	rBV	996486	1512678	1.31%	0.794%
18	13.904	1819	1822	1830	rVB5	99331	211957	0.18%	0.111%
19	14.021	1837	1842	1848	rBV2	244319	467701	0.40%	0.246%
20	14.110	1853	1857	1863	rVB	258989	397045	0.34%	0.208%
21	14.298	1885	1889	1897	rBV3	208930	432872	0.37%	0.227%
22	14.410	1904	1908	1915	rBV4	266940	417973	0.36%	0.219%
23	14.480	1916	1920	1926	rBV2	684562	1020959	0.88%	0.536%
24	14.598	1935	1940	1946	rBV	2185320	3357379	2.90%	1.763%
25	14.657	1947	1950	1957	rVB7	202197	299851	0.26%	0.157%
26	15.151	2030	2034	2038	rBV	592035	935010	0.81%	0.491%
27	15.233	2038	2048	2056	rVV	2739161	5281846	4.56%	2.773%
28	15.421	2076	2080	2086	rBV2	1347563	2307870	1.99%	1.212%
29	15.539	2096	2100	2106	rBV	610755	947713	0.82%	0.498%
30	15.598	2106	2110	2114	rBV2	334167	563151	0.49%	0.296%
31	15.704	2123	2128	2135	rVB	2504100	3733378	3.22%	1.960%
32	17.033	2344	2354	2356	rBV	45018509	115868452	100.00%	60.835%
33	17.145	2369	2373	2378	rBV4	2518435	4752303	4.10%	2.495%
34	20.621	2962	2964	2968	rVB2	2684307	2866749	2.47%	1.505%
35	21.450	3101	3105	3113	rVB	10834491	13355122	11.53%	7.012%
36	21.544	3117	3121	3130	rVB2	3480818	5915315	5.11%	3.106%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

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-BM121523.MA.M
Title : SVOA CALIBRATION

37	23.968	3529	3533	3541	rVB	1951187	4030514	3.48%	2.116%
38	25.709	3820	3829	3845	rVB	4972768	13644185	11.78%	7.164%

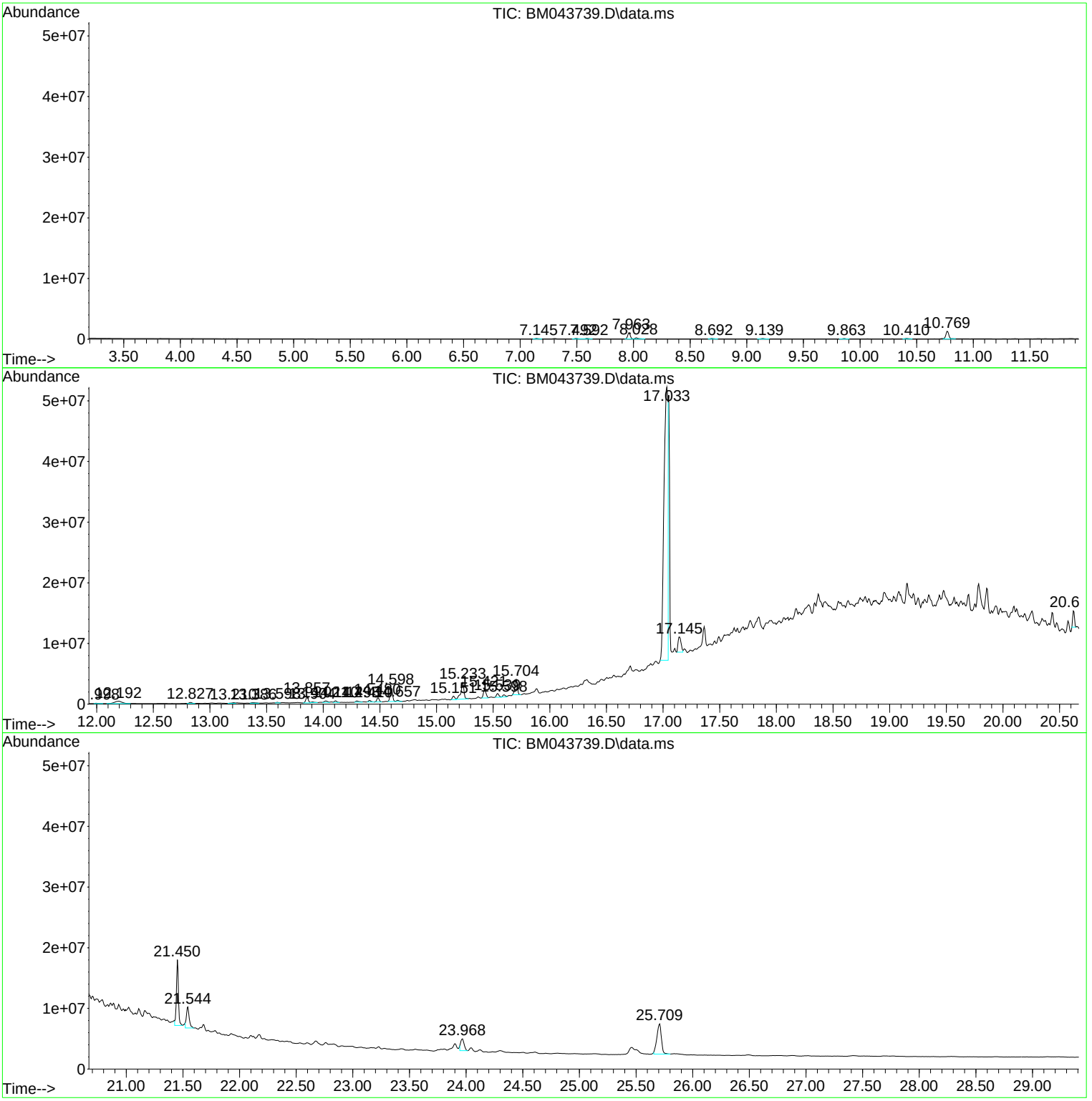
Sum of corrected areas: 190462010

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

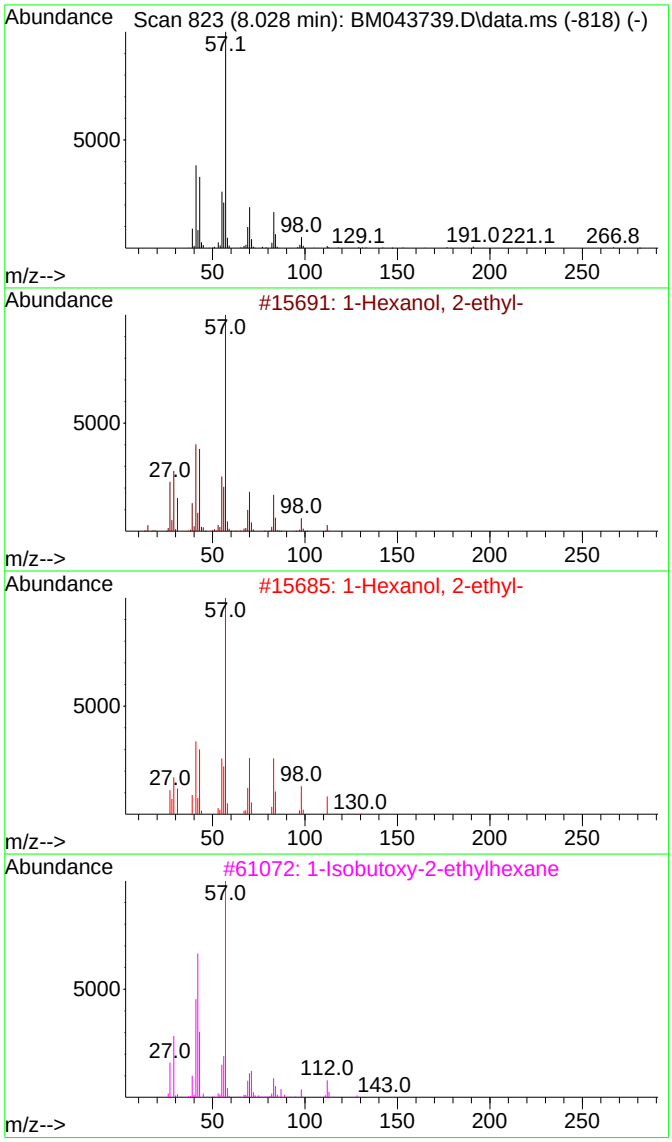
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

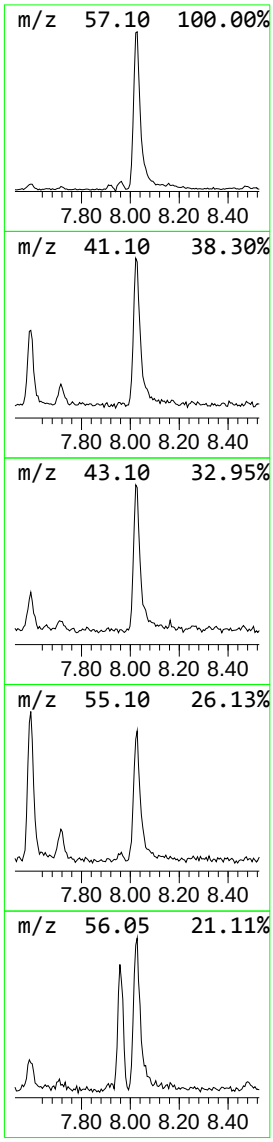
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.028	3.90 ng/ul	324599	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
3	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	72
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

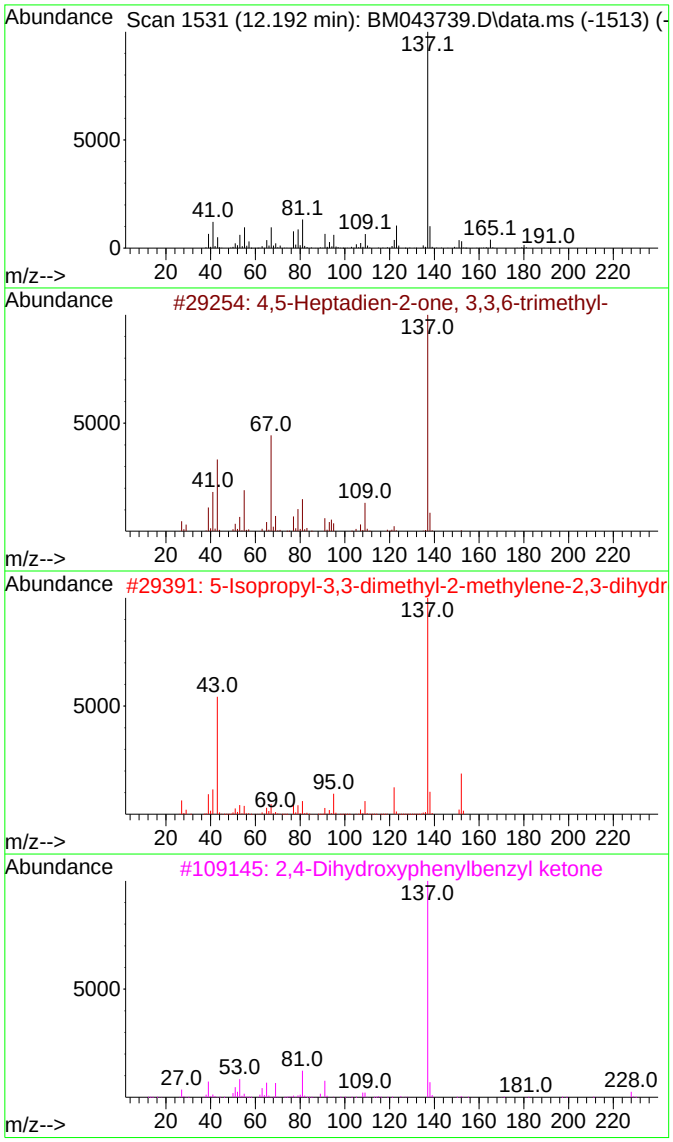
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

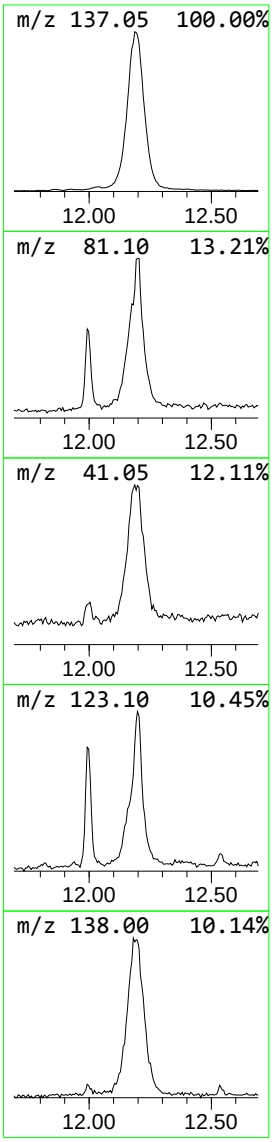
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 4,5-Heptadien-2-one, 3,3,6-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.192	16.89 ng/ul	1865440	Naphthalene-d8	10.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5-Heptadien-2-one, 3,3,6-trime...	152	C10H16O	081250-41-1	50
2	5-Isopropyl-3,3-dimethyl-2-methy...	152	C10H16O	081250-44-4	50
3	2,4-Dihydroxyphenylbenzyl ketone	228	C14H12O3	003669-41-8	42
4	Guaiacol, 4-butyl-	180	C11H16O2	059832-96-1	40
5	2-Pentenoic acid, 3-methyl-5-(2,...	236	C15H24O2	1000196-81-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

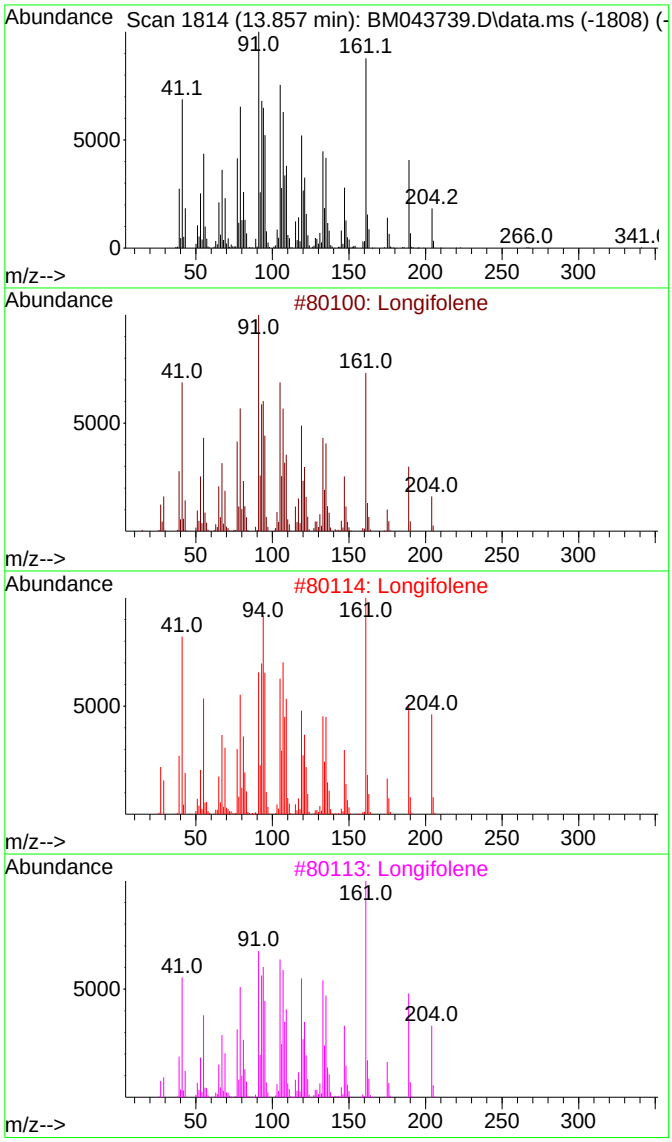
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

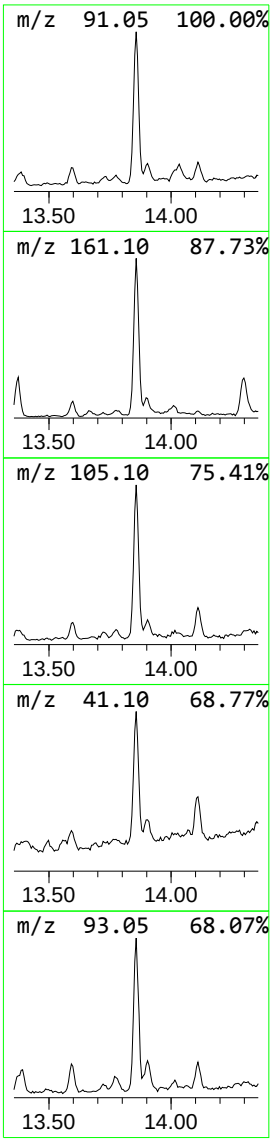
****

Peak Number 3 Longifolene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	9.01 ng/ul	1512680	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Longifolene	204	C15H24	000475-20-7	99
2	Longifolene	204	C15H24	000475-20-7	99
3	Longifolene	204	C15H24	000475-20-7	98
4	Longifolene	204	C15H24	000475-20-7	94
5	Longifolene	204	C15H24	000475-20-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

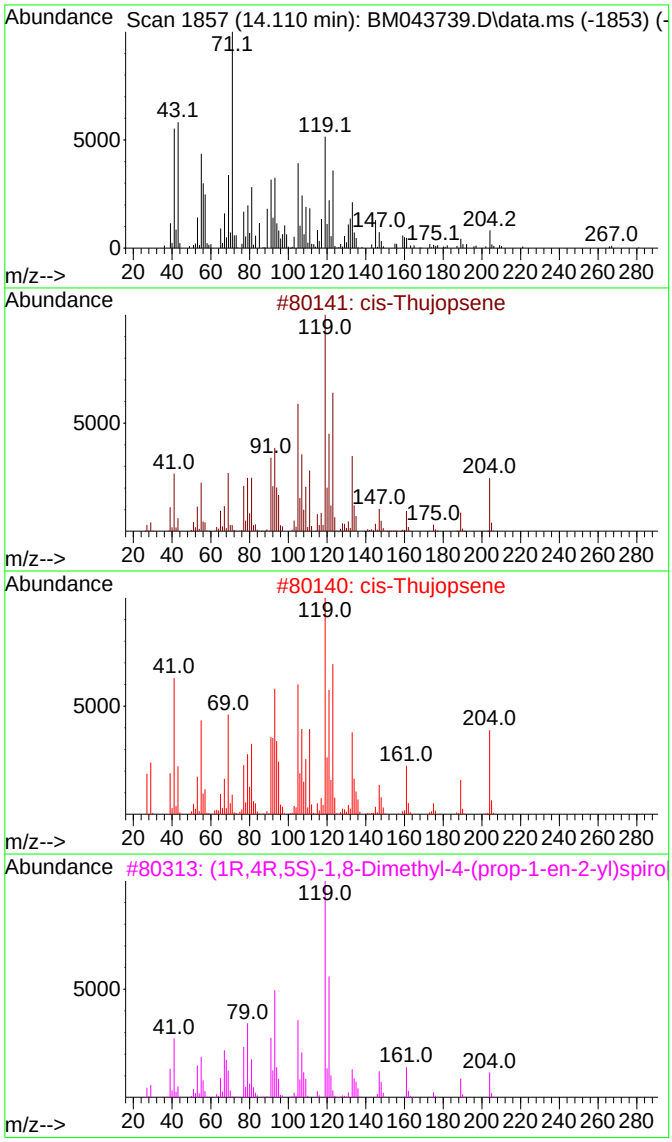
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

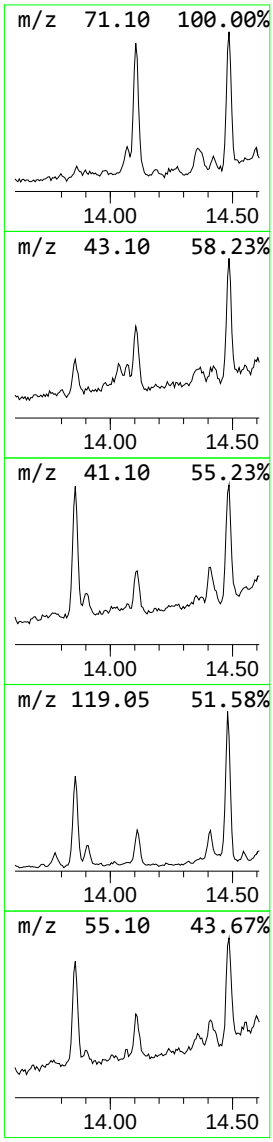
****

Peak Number 4 cis-Thujopsene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.110	2.37 ng/ul	397045	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-Thujopsene	204	C15H24	000470-40-6	95
2	cis-Thujopsene	204	C15H24	000470-40-6	43
3	(1R,4R,5S)-1,8-Dimethyl-4-(prop-...	204	C15H24	729602-94-2	42
4	Tricyclo[5.4.0.0(2,8)]undec-9-en...	204	C15H24	005989-08-2	42
5	Tricyclo[5.4.0.0(2,8)]undec-9-en...	204	C15H24	005989-08-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

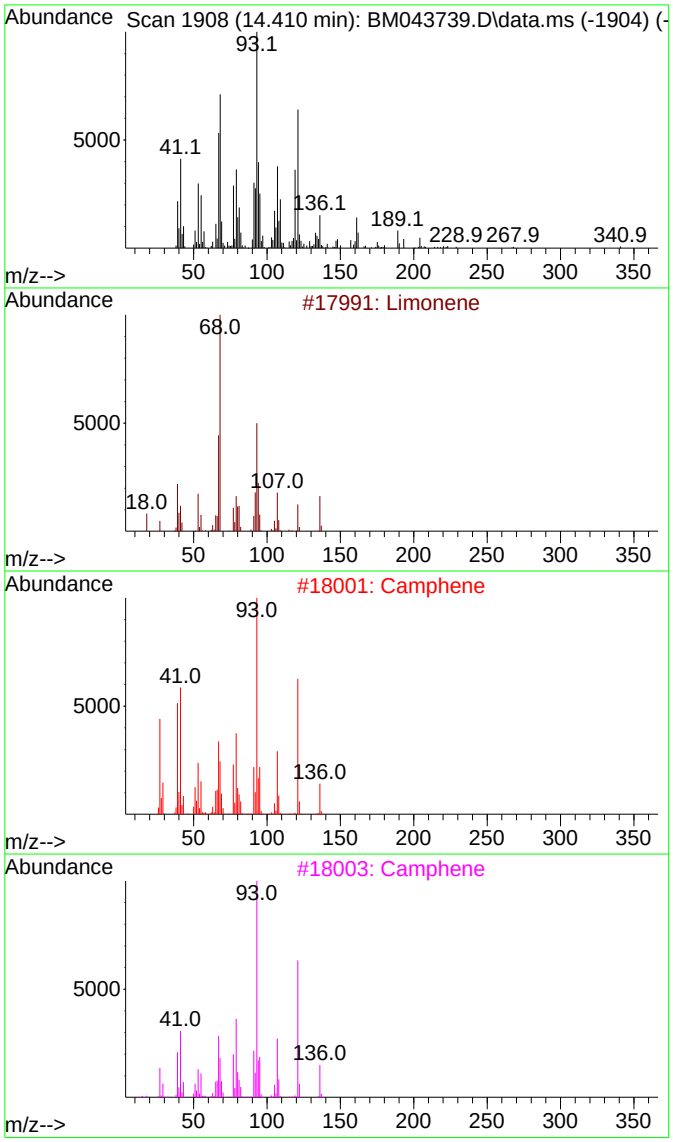
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

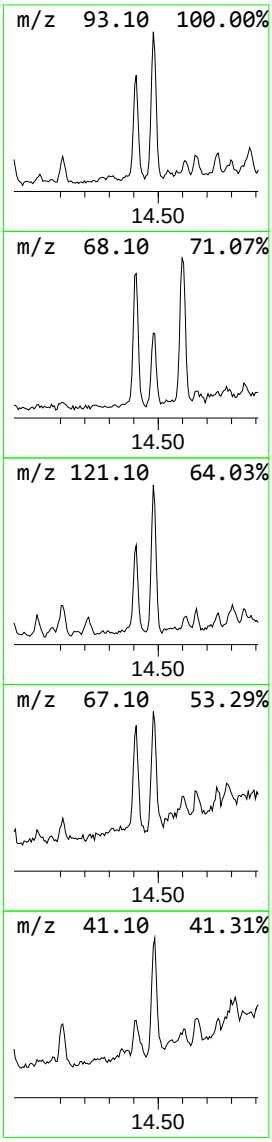
****

Peak Number 5 Limonene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.410	2.49 ng/ul	417973	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Limonene	136	C10H16	000138-86-3	76
2	Camphene	136	C10H16	000079-92-5	55
3	Camphene	136	C10H16	000079-92-5	55
4	1H-Benzocycloheptene, 2,4a,5,6,7...	204	C15H24	003853-83-6	49
5	Camphene	136	C10H16	000079-92-5	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

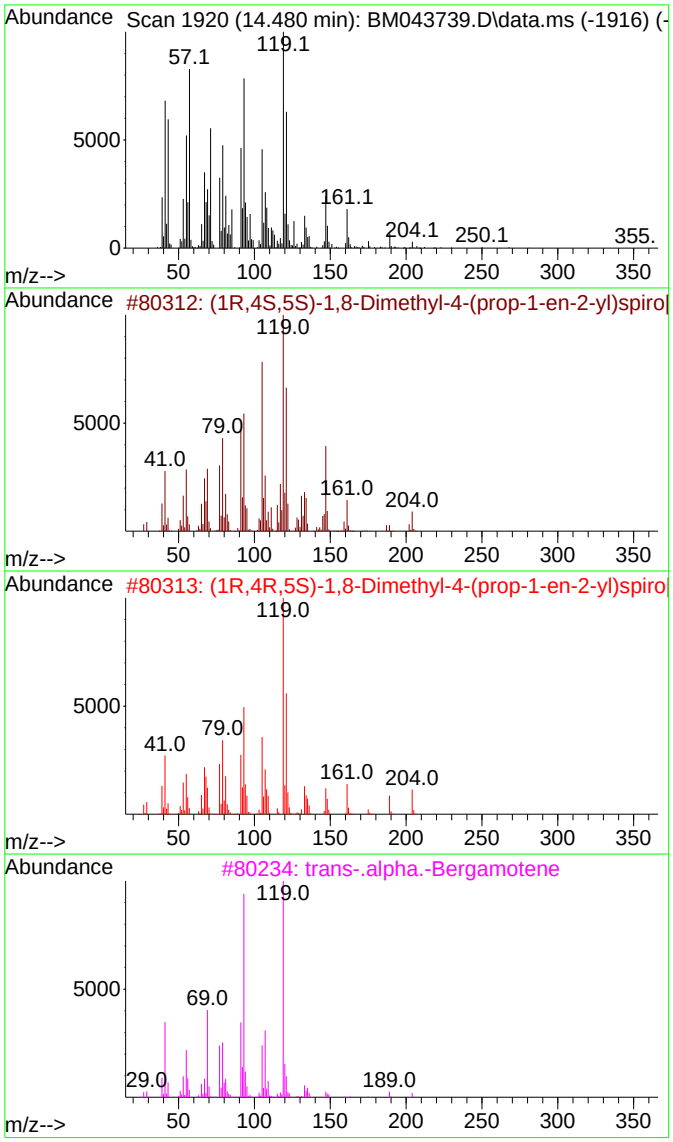
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

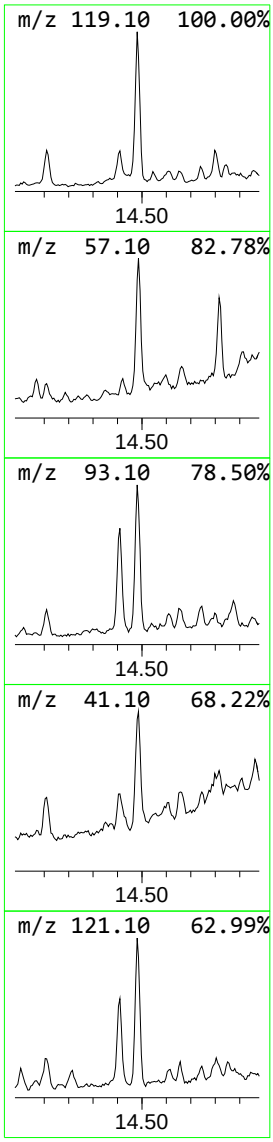
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (1R,4S,5S)-1,8-Dimethyl-4-(... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	6.08 ng/ul	1020960	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R,4S,5S)-1,8-Dimethyl-4-(prop-...	204	C15H24	043219-80-3	83
2	(1R,4R,5S)-1,8-Dimethyl-4-(prop-...	204	C15H24	729602-94-2	74
3	trans-.alpha.-Bergamotene	204	C15H24	013474-59-4	52
4	trans-.alpha.-Bergamotene	204	C15H24	013474-59-4	52
5	cis-.alpha.-Bergamotene	204	C15H24	018252-46-5	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

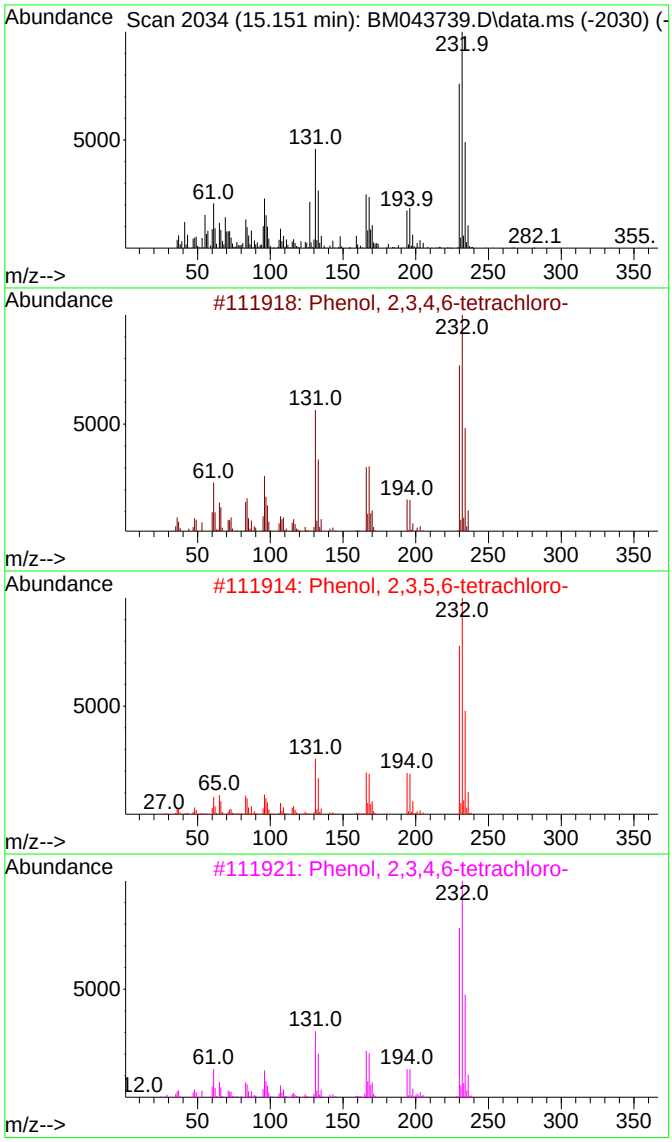
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

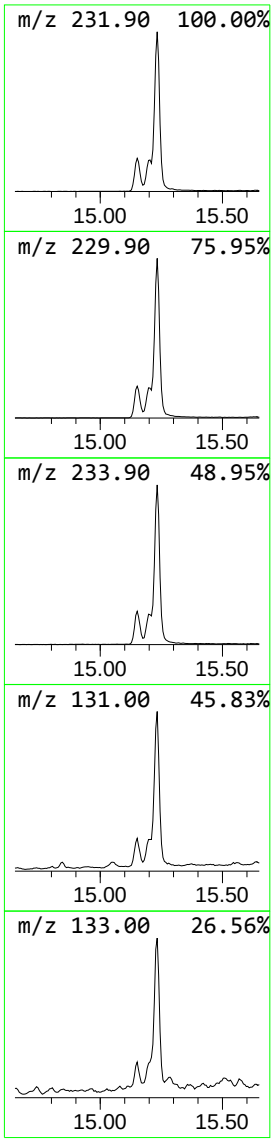
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.151	5.57 ng/ul	935010	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
3	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
4	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
5	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

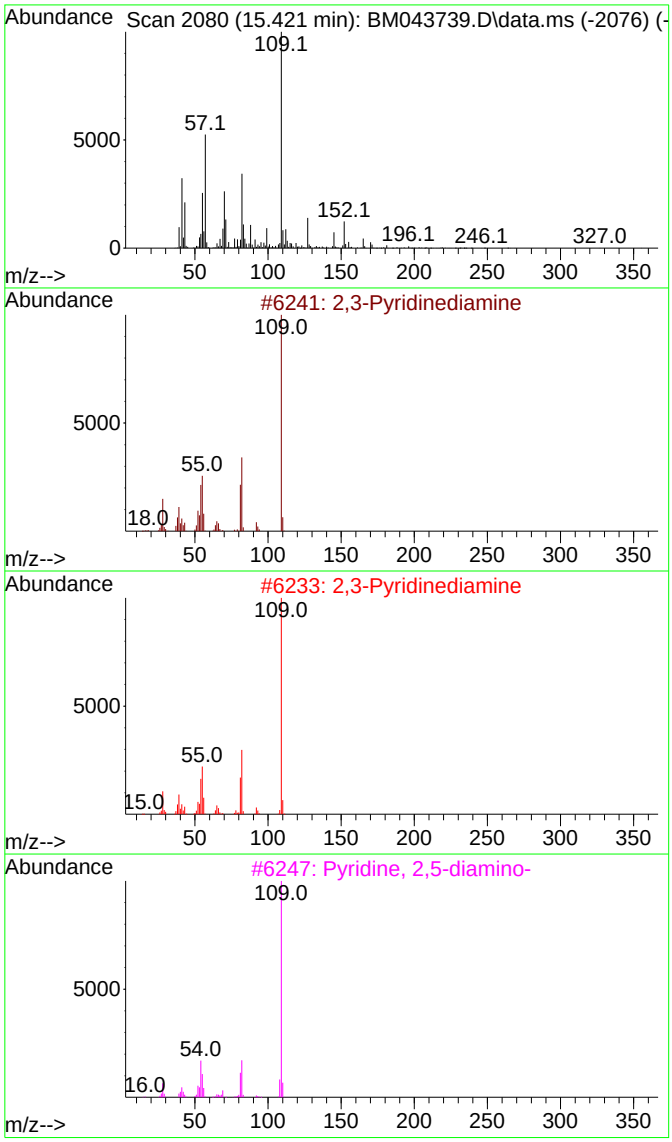
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

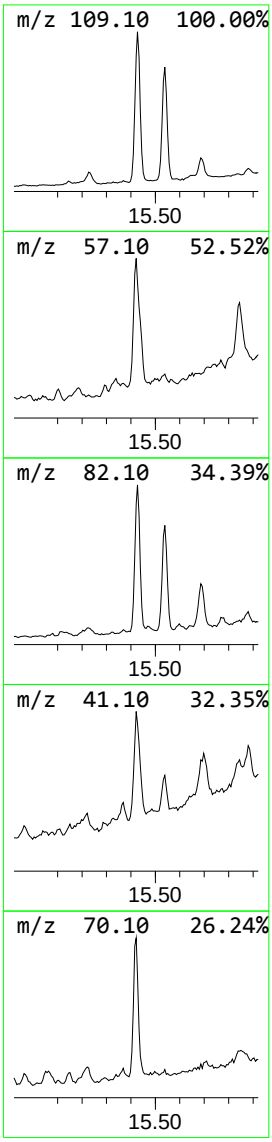
****

Peak Number 8 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.421	13.75 ng/ul	2307870	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Pyridinediamine	109	C5H7N3	000452-58-4	43
2	2,3-Pyridinediamine	109	C5H7N3	000452-58-4	43
3	Pyridine, 2,5-diamino-	109	C5H7N3	004318-76-7	43
4	1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	38
5	Pyridine-2,4-diamine	109	C5H7N3	000461-88-1	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

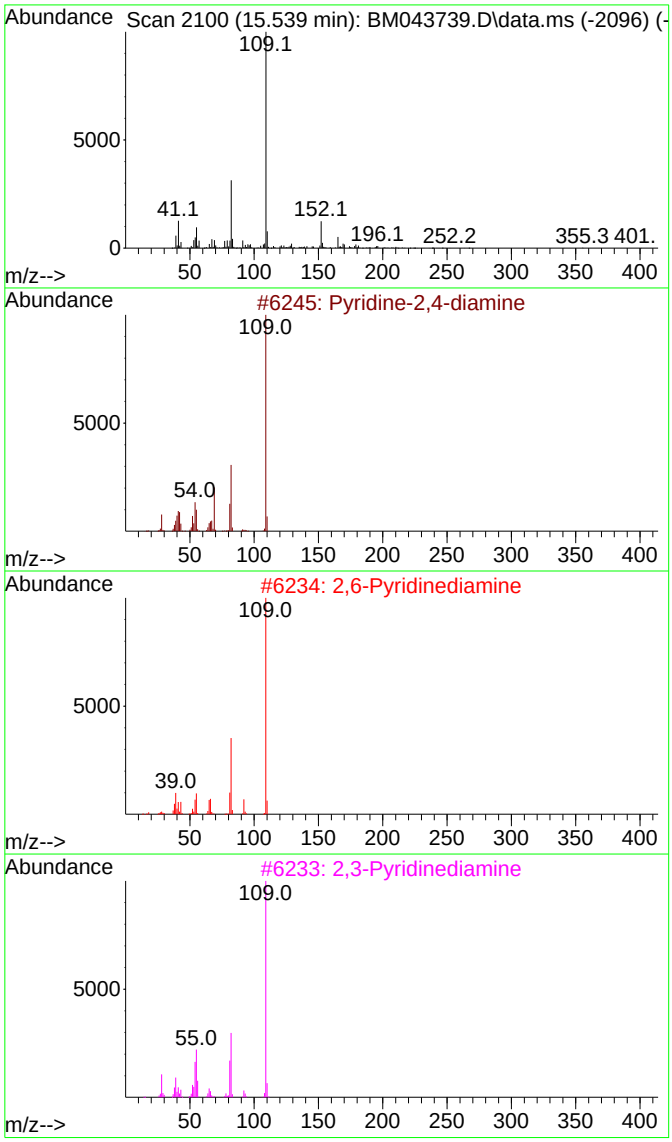
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

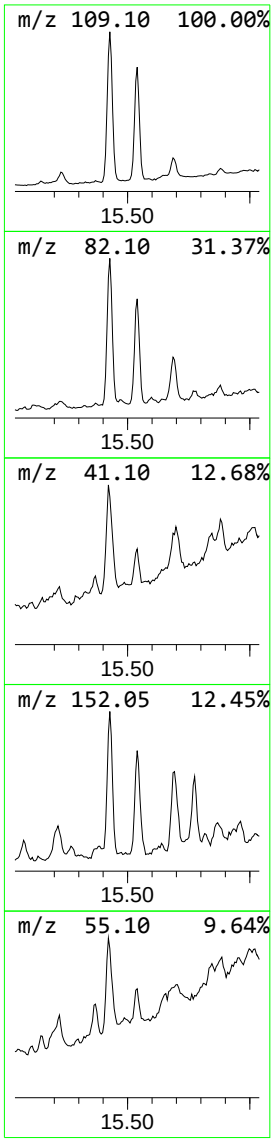
****

Peak Number 9 Pyridine-2,4-diamine Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.539	5.65 ng/ul	947713	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine-2,4-diamine	109	C5H7N3	000461-88-1	64
2	2,6-Pyridinediamine	109	C5H7N3	000141-86-6	64
3	2,3-Pyridinediamine	109	C5H7N3	000452-58-4	50
4	(4-Fluorophenyl) methanol, tert....	182	C11H15FO	1000374-64-1	50
5	(2-Fluorophenyl) methanol, tert....	182	C11H15FO	1000374-56-8	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

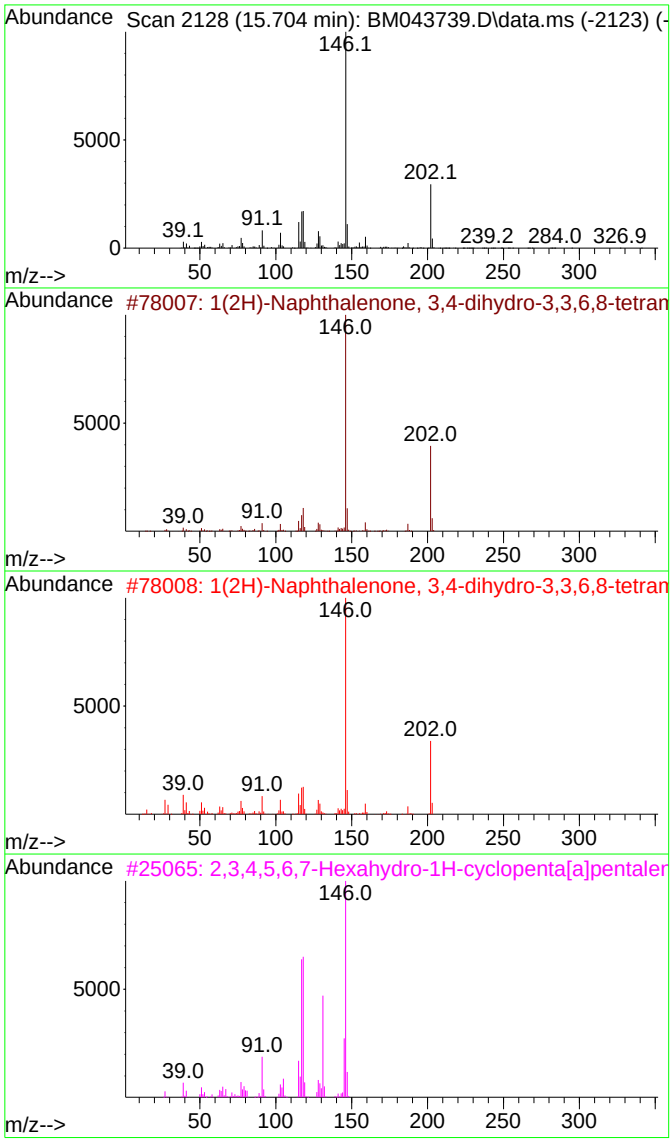
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

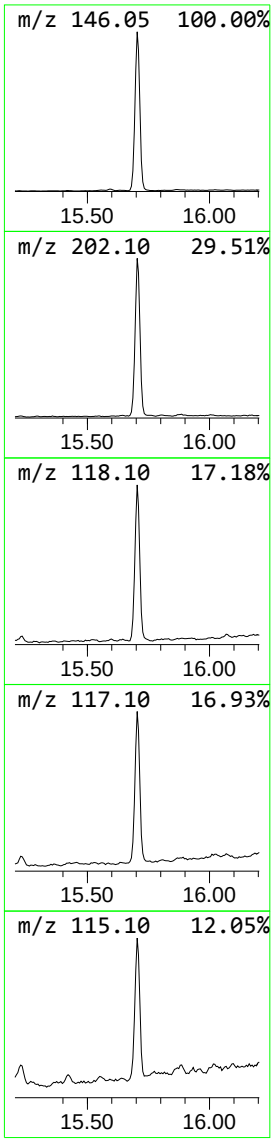
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.704	22.24 ng/ul	3733380	Acenaphthene-d10	14.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4-dihydro...	202	C14H18O	005409-55-2	97
2	1(2H)-Naphthalenone, 3,4-dihydro...	202	C14H18O	005409-55-2	94
3	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	53
4	Phthalazin-1-one	146	C8H6N2O	000119-39-1	50
5	4-Pyrimidinol, 5-fluoro-6-(fluor...	146	C5H4F2N2O	1010319-28-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

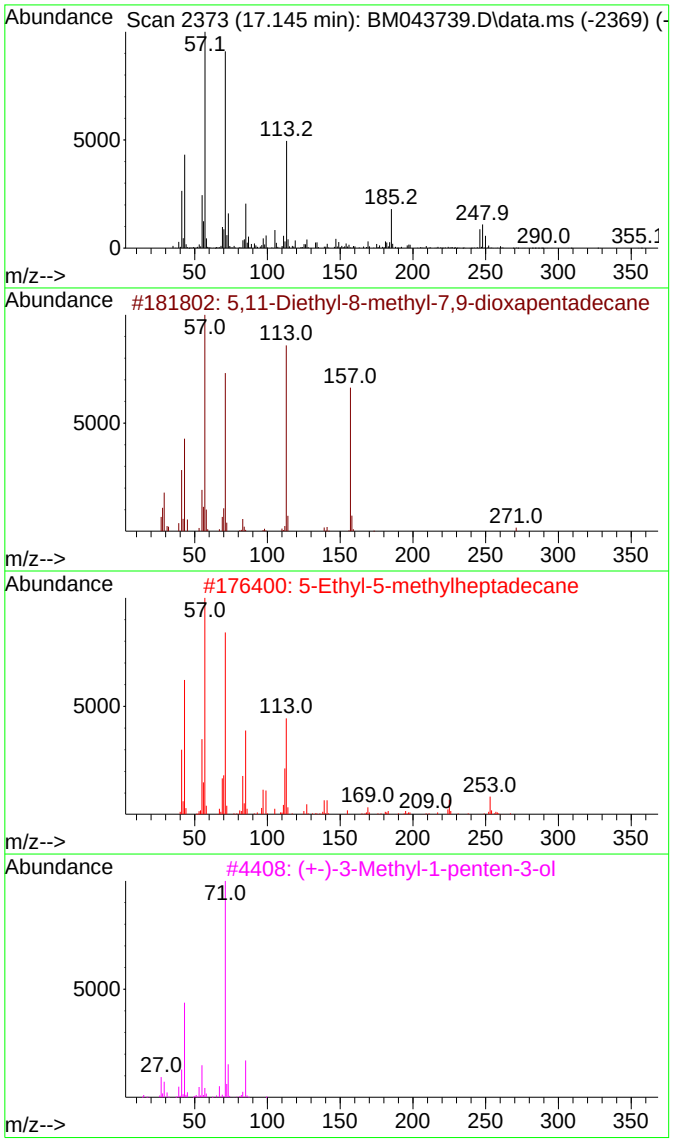
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

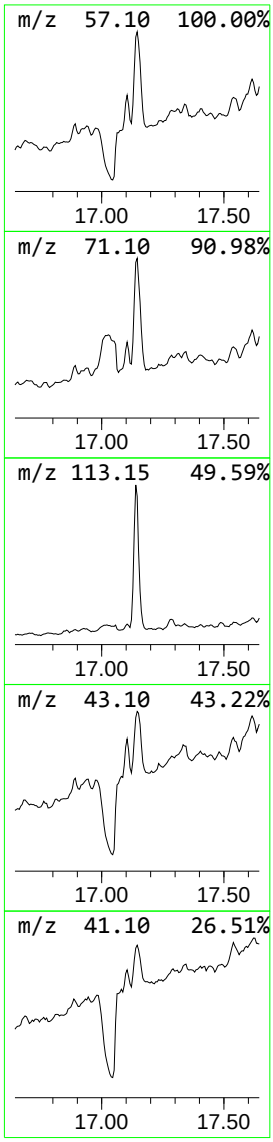
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 5,11-Diethyl-8-methyl-7,9-d... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.145	20.00 ng/ul	4752300	Phenanthrene-d10	17.362

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,11-Diethyl-8-methyl-7,9-dioxap...	286	C18H38O2	1000334-83-1	53
2	5-Ethyl-5-methylheptadecane	282	C20H42	1000360-42-5	53
3	(+-)-3-Methyl-1-penten-3-ol	100	C6H12O	1010238-75-4	46
4	Octane, 1-iodo-	240	C8H17I	000629-27-6	43
5	Octane, 1-iodo-	240	C8H17I	000629-27-6	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

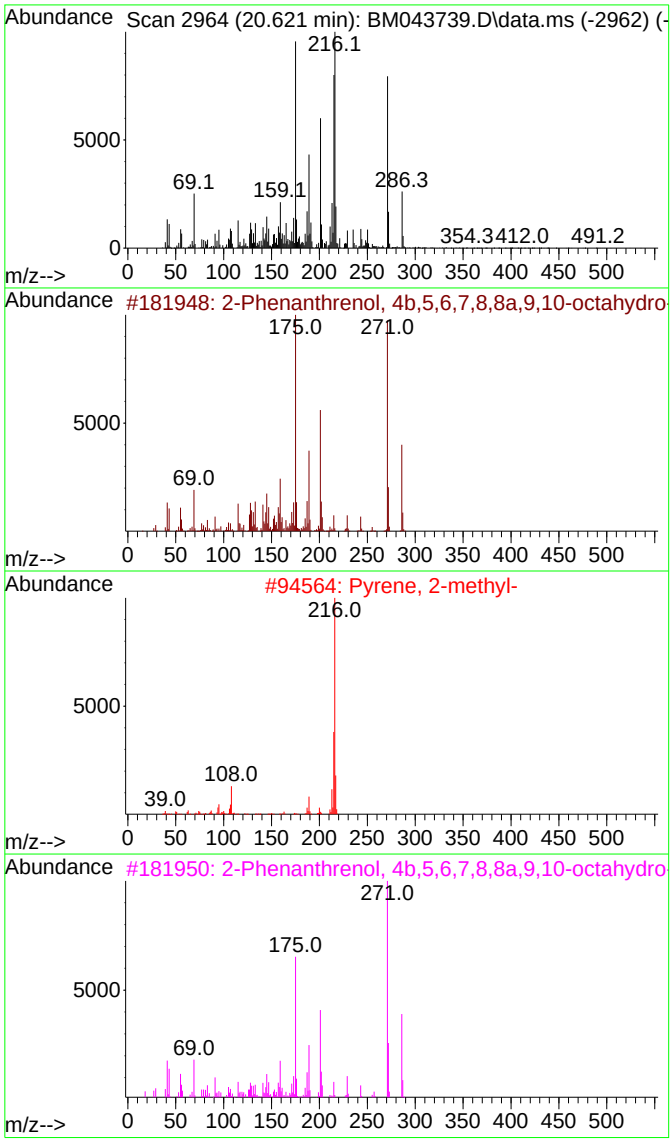
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

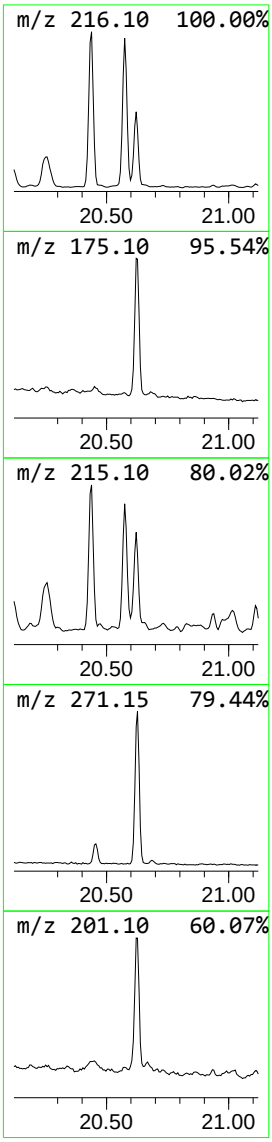
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.621	9.69 ng/ul	2866750	Chrysene-d12	21.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	97
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	56
3	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	55
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	46
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

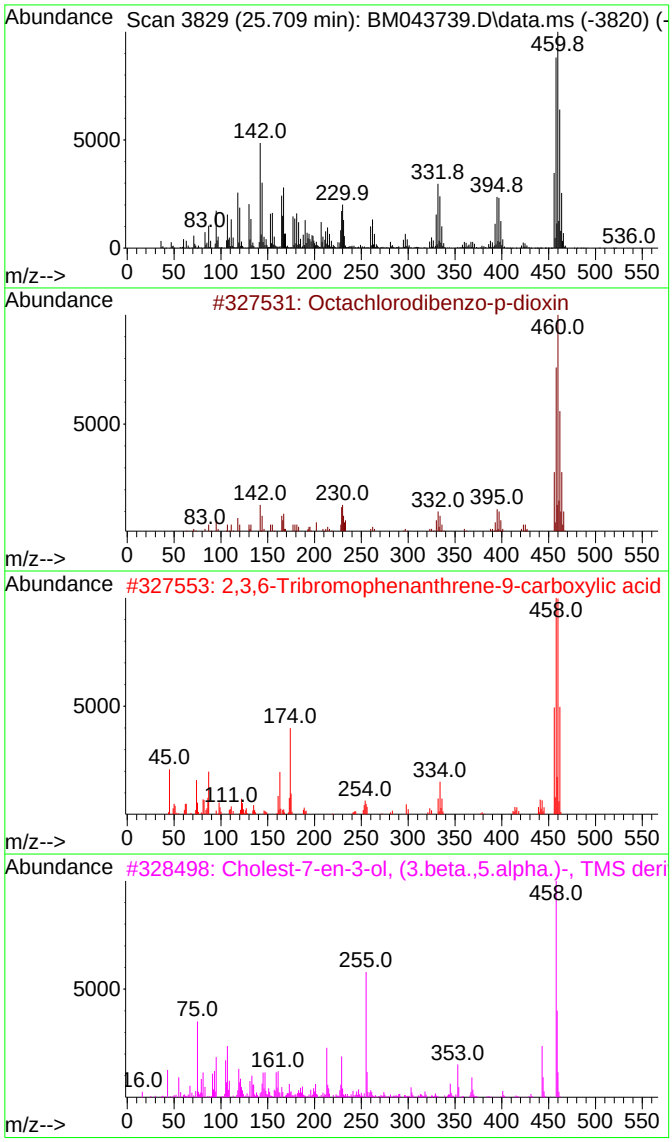
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

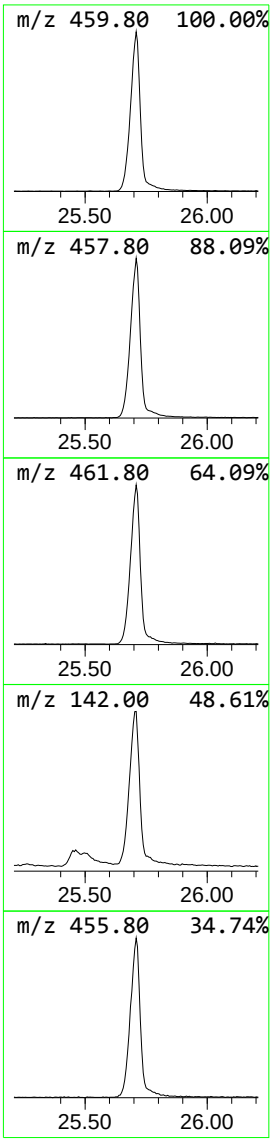
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Octachlorodibenzo-p-dioxin Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.709	67.70 ng/ul	13644200	Perylene-d12	23.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	87
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	43
3	Cholest-7-en-3-ol, (3.beta.,5.al...	458	C30H54OSi	002665-03-4	11
4	Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10
5	2'-Hydroxy-3,3',4'-trimethoxycha...	460	C21H17F5O6	1000453-84-7	9





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
Data File : BM043739.D
Acq On : 03 Jan 2024 18:44
Operator : MA/JU
Sample : 05952-08DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF91DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.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
1-Hexanol, 2-et...	8.028	3.9	ng/ul	324599	1	7.963	1665610	20.0
4,5-Heptadien-2...	12.192	16.9	ng/ul	1865440	2	10.769	2209320	20.0
Longifolene	13.857	9.0	ng/ul	1512680	3	14.598	3357380	20.0
cis-Thujopsene	14.110	2.4	ng/ul	397045	3	14.598	3357380	20.0
Limonene	14.410	2.5	ng/ul	417973	3	14.598	3357380	20.0
(1R,4S,5S)-1,8-...	14.480	6.1	ng/ul	1020960	3	14.598	3357380	20.0
Phenol, 2,3,5,6...	15.151	5.6	ng/ul	935010	3	14.598	3357380	20.0
unknown-01	15.421	13.8	ng/ul	2307870	3	14.598	3357380	20.0
Pyridine-2,4-di...	15.539	5.7	ng/ul	947713	3	14.598	3357380	20.0
1(2H)-Naphthale...	15.704	22.2	ng/ul	3733380	3	14.598	3357380	20.0
5,11-Diethyl-8-...	17.145	20.0	ng/ul	4752300	4	17.362	4752300	20.0
2-Phenanthrenol...	20.621	9.7	ng/ul	2866750	5	21.539	5915320	20.0
Octachlorodiben...	25.709	67.7	ng/ul	13644200	6	23.968	4030510	20.0