LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

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_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP017512.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.834	122	127	135	rVB	88426	117238	0.17%	0.041%
2	3.911	135	140	150	rVB	66866	100288	0.15%	0.035%
3	7.093	675	681	697	rBV2	361399	823541	1.21%	0.288%
4	7.275	705	712	722	rBV	329534	511641	0.75%	0.179%
5	7.452	735	742	755	rBV	356386	593717	0.87%	0.207%
6	7.934	818	824	836	rBV	458343	720771	1.06%	0.252%
7	8.657	937	947	955	rBV	281043	470732	0.69%	0.164%
8	8.793	964	970	978	rVB	122083	201770	0.30%	0.071%
9	9.140	1023	1029	1039	rVB	380115	624407	0.92%	0.218%
10	9.857	1145	1151	1158	rBV	211761	351321	0.52%	0.123%
11	10.216	1203	1212	1224	rBV5	30619	100384	0.15%	0.035%
12	10.393	1235	1242	1254	rVB	340788	600653	0.88%	0.210%
13	10.681	1284	1291	1299	rVB	110767	189683	0.28%	0.066%
14	10.775	1299	1307	1313	rVV	578411	941929	1.39%	0.329%
15	10.946	1328	1336	1344	rBV	192124	361934	0.53%	0.126%
16	11.122	1360	1366	1374	rBV	70836	126445	0.19%	0.044%
17	12.116	1526	1535	1546	rBV	918523	1493785	2.20%	0.522%
18	13.434	1754	1759	1763	rBV	127667	193936	0.29%	0.068%
19	13.810	1819	1823	1826	rBV	103825	138715	0.20%	0.048%
20	13.851	1826	1830	1835	rVV	176879	239478	0.35%	0.084%
21	14.045	1857	1863	1869	rVV	819798	1074805	1.58%	0.376%
22	14.328	1903	1911	1918	rBV	947296	1498106	2.21%	0.523%
23	14.404	1920	1924	1929	rVB3	97102	148864	0.22%	0.052%
24	14.634	1957	1963	1968	rVV	785214	1162307	1.71%	0.406%
25	14.898	2004	2008	2018	rBV5	50026	153524	0.23%	0.054%
26	15.375	2085	2089	2095	rBV2	121219	182793	0.27%	0.064%
27	15.634	2128	2133	2141	rBV	1319498	1869571	2.75%	0.653%
28	16.304	2244	2247	2252	rBV2	100202	158992	0.23%	0.056%
29	17.422	2433	2437	2442	rBV	1002915	1347938	1.98%	0.471%
30	17.522	2450	2454	2461	rBV2	1258923	1831154	2.70%	0.640%
31	19.780	2834	2838	2843	rBV	2037990	2865848	4.22%	1.001%
32	22.104	3228	3233	3244	rVB	8288802	12038168	17.73%	4.206%
33	22.680	3324	3331	3340	rVB	17471336	30671558	45.17%	10.717%
34	22.827	3352	3356	3359	rBV	2429541	3887688	5.72%	1.358%
35	22.863	3359	3362	3366	rVB	3674783	4803025	7.07%	1.678%
36	23.039	3387	3392	3403	rVB2	7693256	16196252	23.85%	5.659%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

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_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

37	23.292	3431	3435	3444	rVB3	2245196	4403999	6.49%	1.539%
38	23.427	3453	3458	3463	rVB	4556240	8803897	12.96%	3.076%
39	23.498	3466	3470	3475	rVB2	3676373	6259268	9.22%	2.187%
40	23.680	3493	3501	3508	rBV	15361249	31625314	46.57%	11.050%
41	23.974	3547	3551	3558	rBV4	1621349	3655874	5.38%	1.277%
42	24.539	3642	3647	3657	rVB4	2532279	6515958	9.60%	2.277%
43	24.704	3671	3675	3683	rVB	2368694	5184618	7.63%	1.812%
44	24.880	3693	3705	3712	rBV	13082590	39351826	57.95%	13.750%
45	25.515	3807	3813	3829	rVB2	3246650	8724205	12.85%	3.048%
46	25.839	3863	3868	3886	rVB2	1956620	6022704	8.87%	2.104%
47	26.780	4021	4028	4042	rVB	2972641	8943081	13.17%	3.125%
48	28.486	4296	4318	4338	rVB	14640642	67908579	100.00%	23.728%

Sum of corrected areas: 286192284

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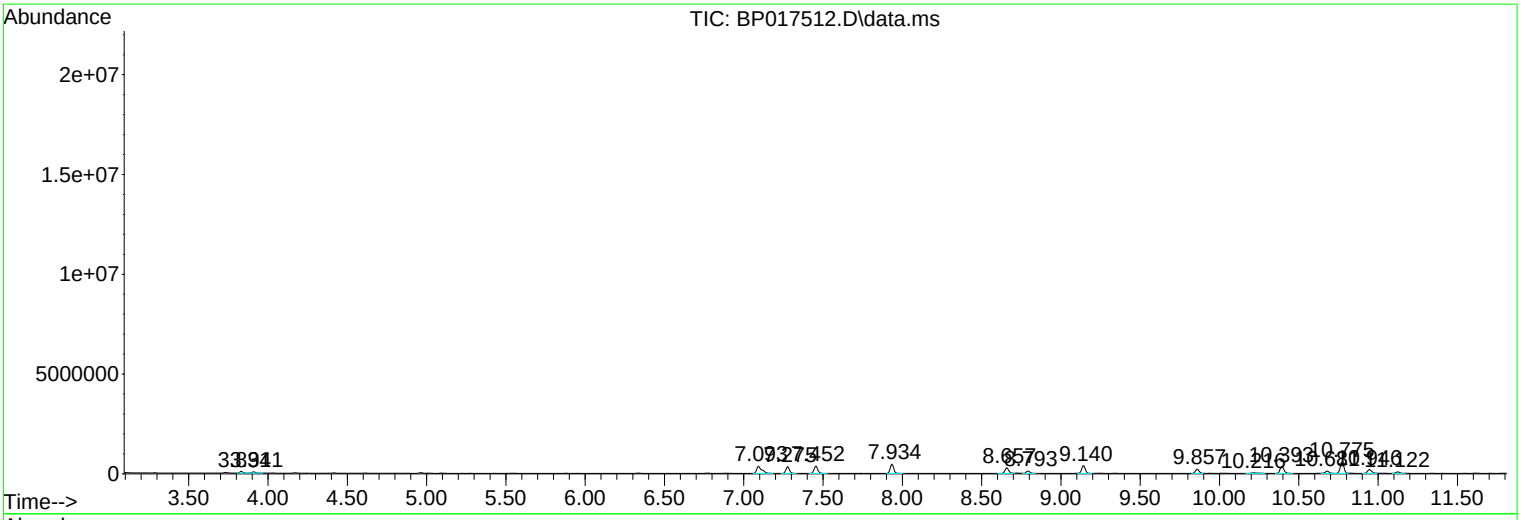
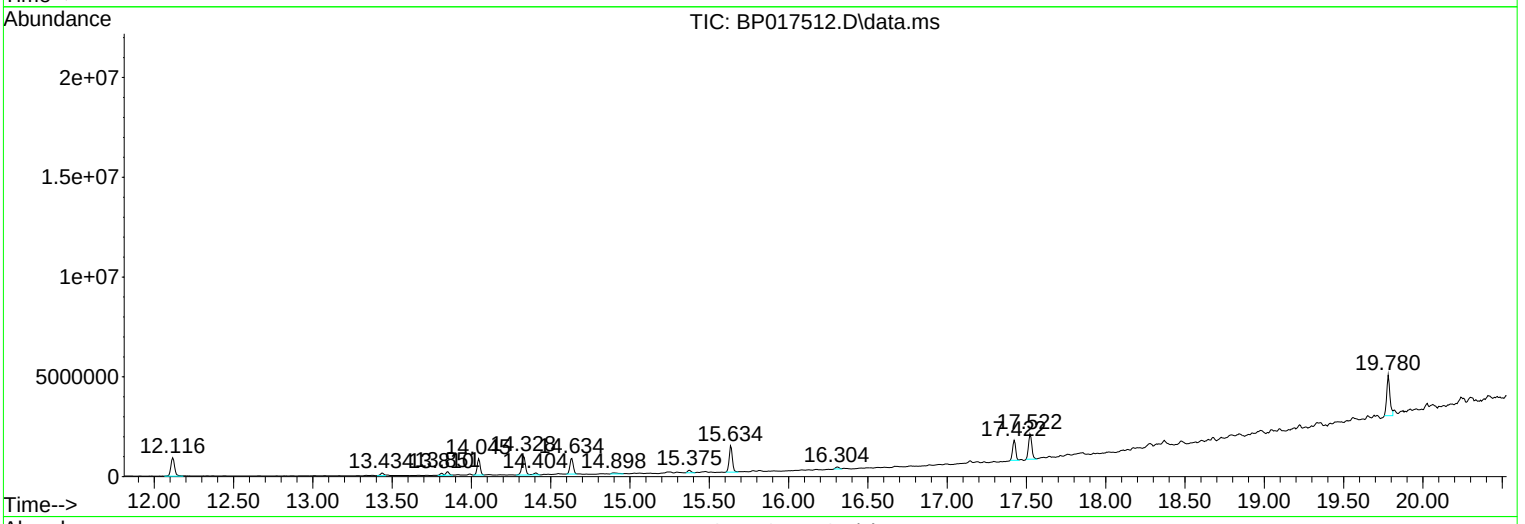
ALS Vial : 36 Sample Multiplier: 1

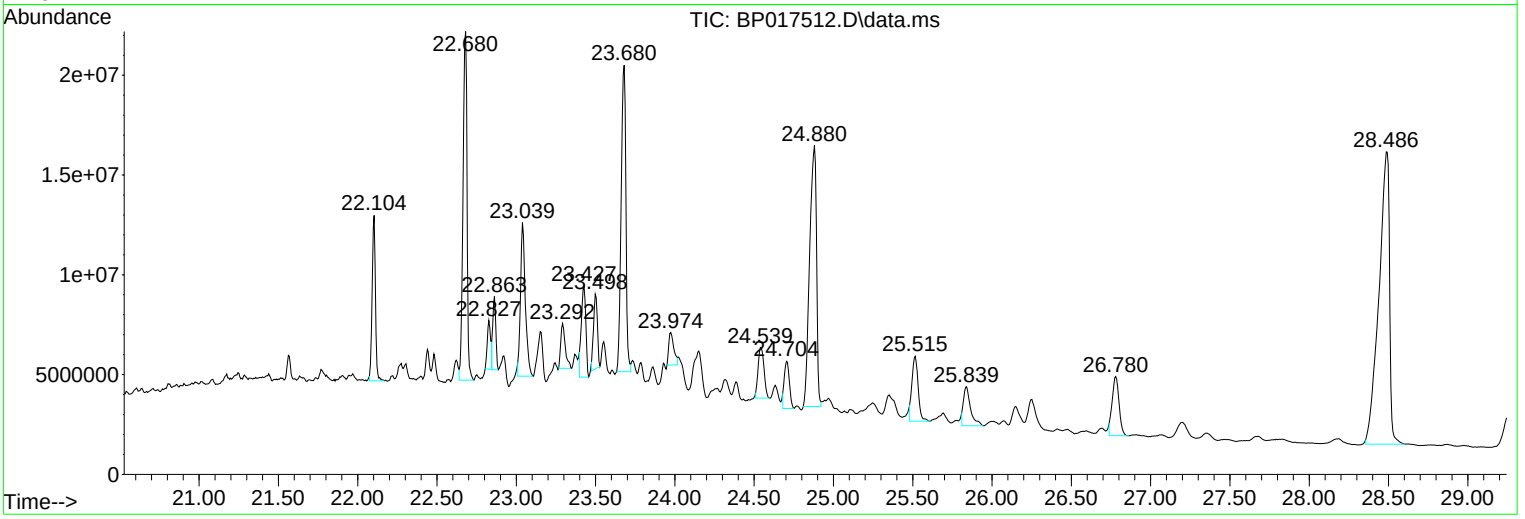
Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

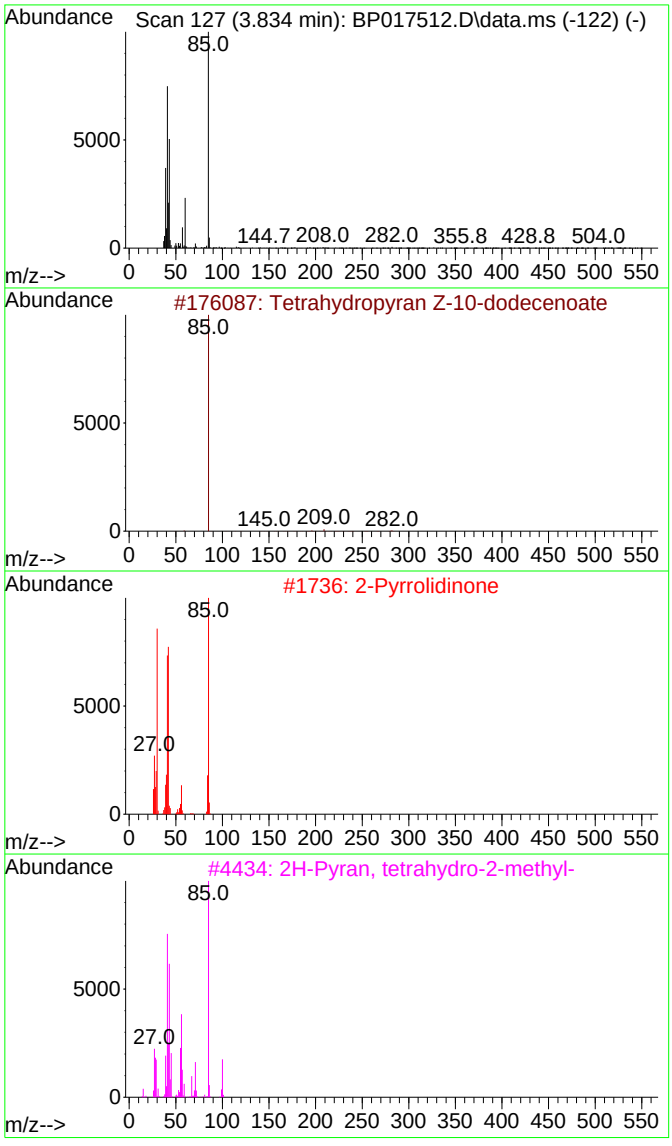
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

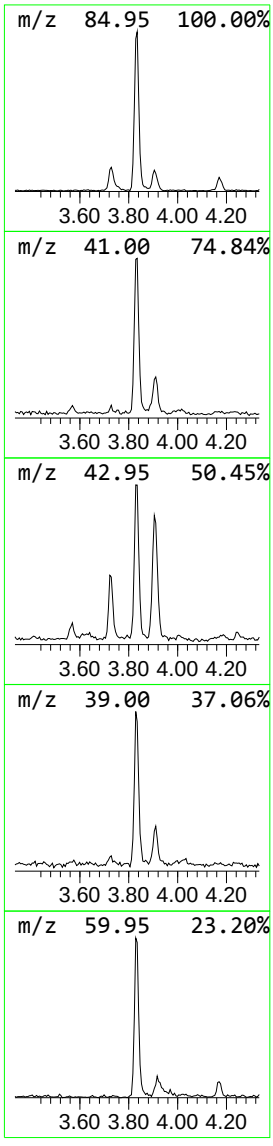
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Tetrahydropyran Z-10-dodece... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.834	3.25 ng/ul	117238	1,4-Dichlorobenzene-d4	7.934

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrahydropyran Z-10-dodecenoate	282	C17H30O3	1000130-96-5	64
2	2-Pyrrolidinone	85	C4H7NO	000616-45-5	40
3	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	39
4	Methacrylamide	85	C4H7NO	000079-39-0	38
5	2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

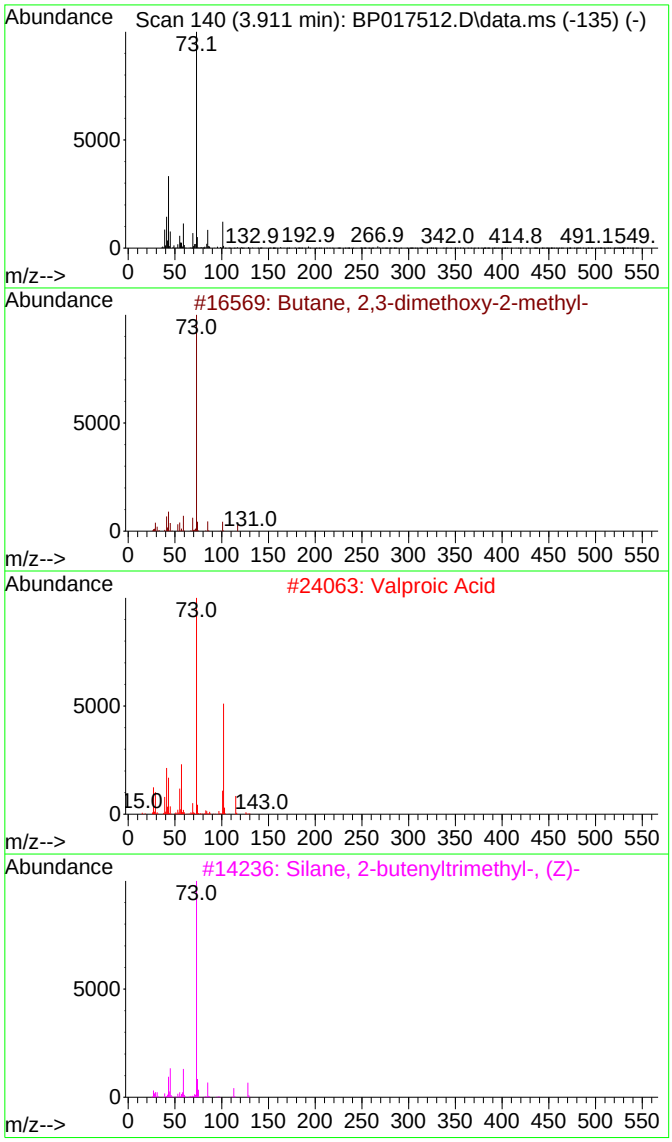
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

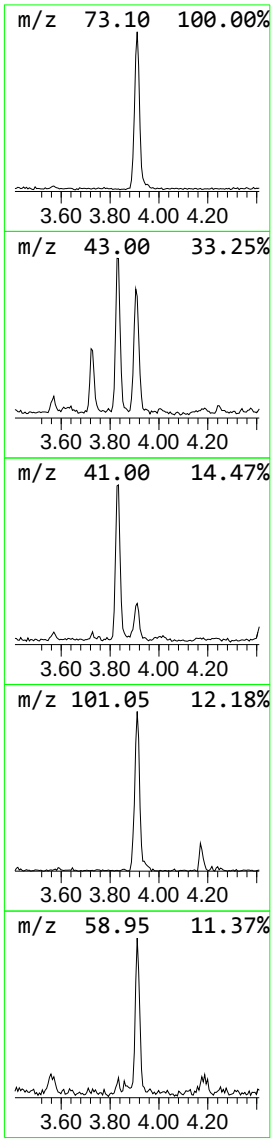
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dimethoxy-2-met... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.911	2.78 ng/ul	100288	1,4-Dichlorobenzene-d4	7.934

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethoxy-2-methyl-	132	C7H16O2	074421-00-4	53
2	Valproic Acid	144	C8H16O2	000099-66-1	38
3	Silane, 2-butenyltrimethyl-, (Z)-	128	C7H16Si	017486-13-4	37
4	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	37
5	3-Heptanol, 3-methyl-	130	C8H18O	005582-82-1	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

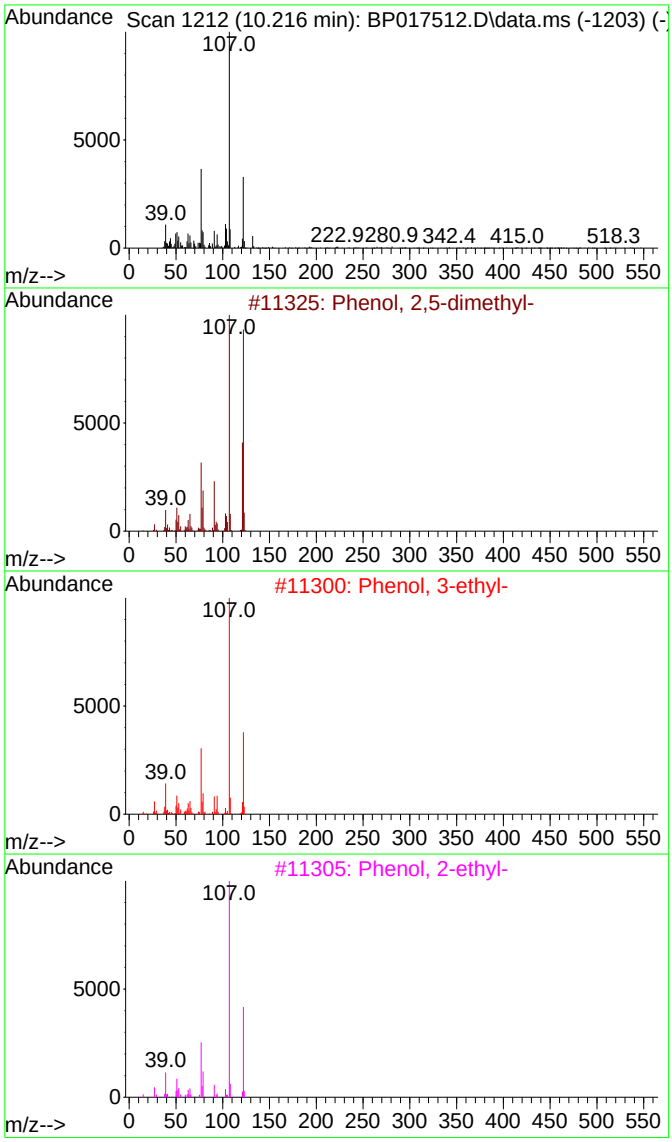
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

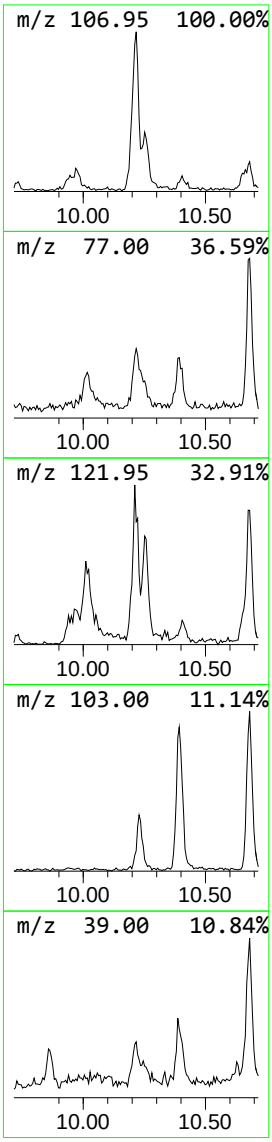
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, 2,5-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.216	2.13 ng/ul	100384	Naphthalene-d8	10.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	86
2	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	81
3	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	74
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	74
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

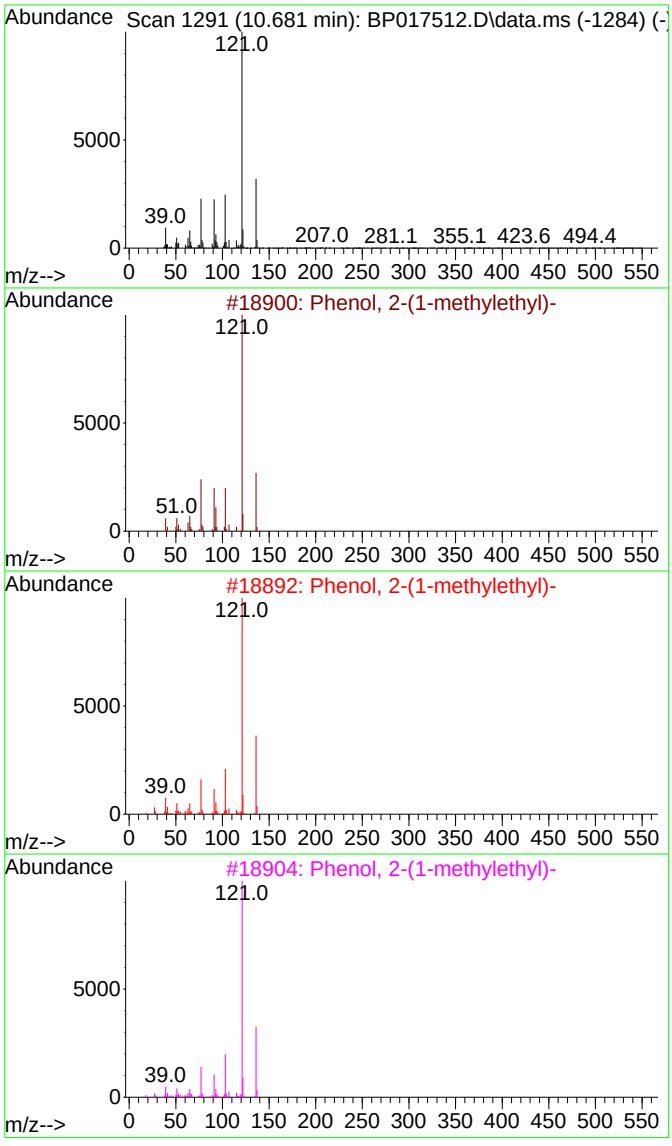
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

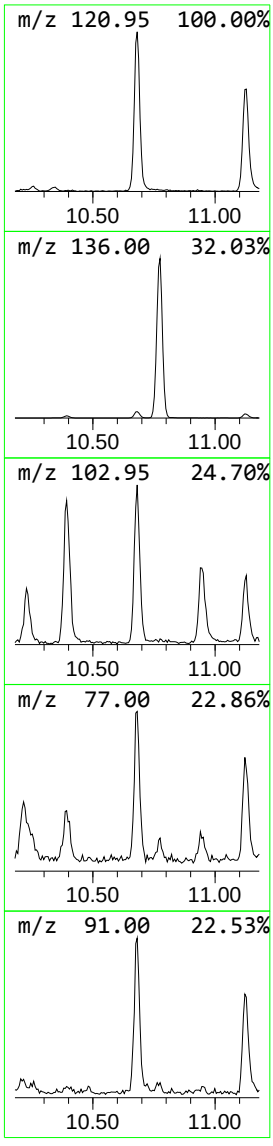
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 2-(1-methylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.681	4.03 ng/ul	189683	Naphthalene-d8	10.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	94
2	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
3	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	90
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

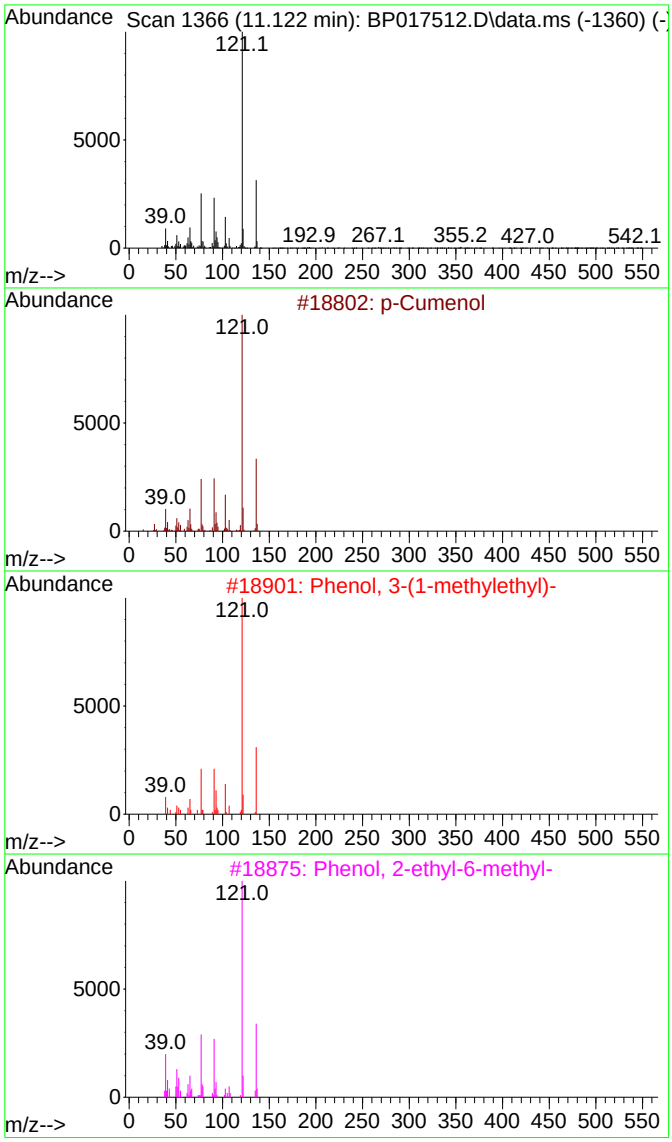
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

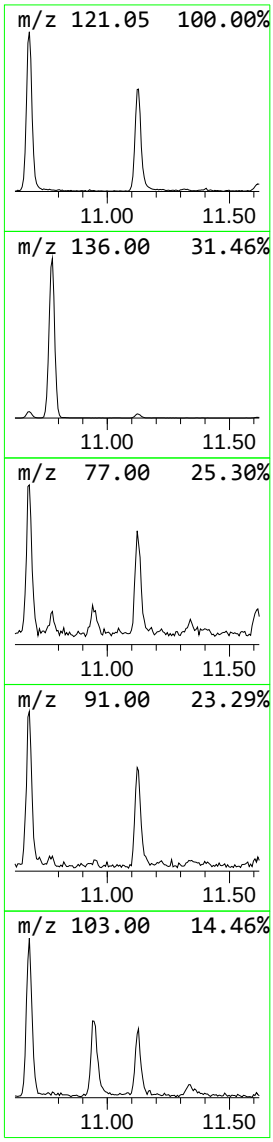
****

Peak Number 5 p-Cumenol Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.122	2.68 ng/ul	126445	Naphthalene-d8	10.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cumenol	136	C9H12O	000099-89-8	94
2	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
3	Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	90
4	3-Isopropylphenyl N-methylcarbam...	193	C11H15NO2	1000470-98-9	90
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

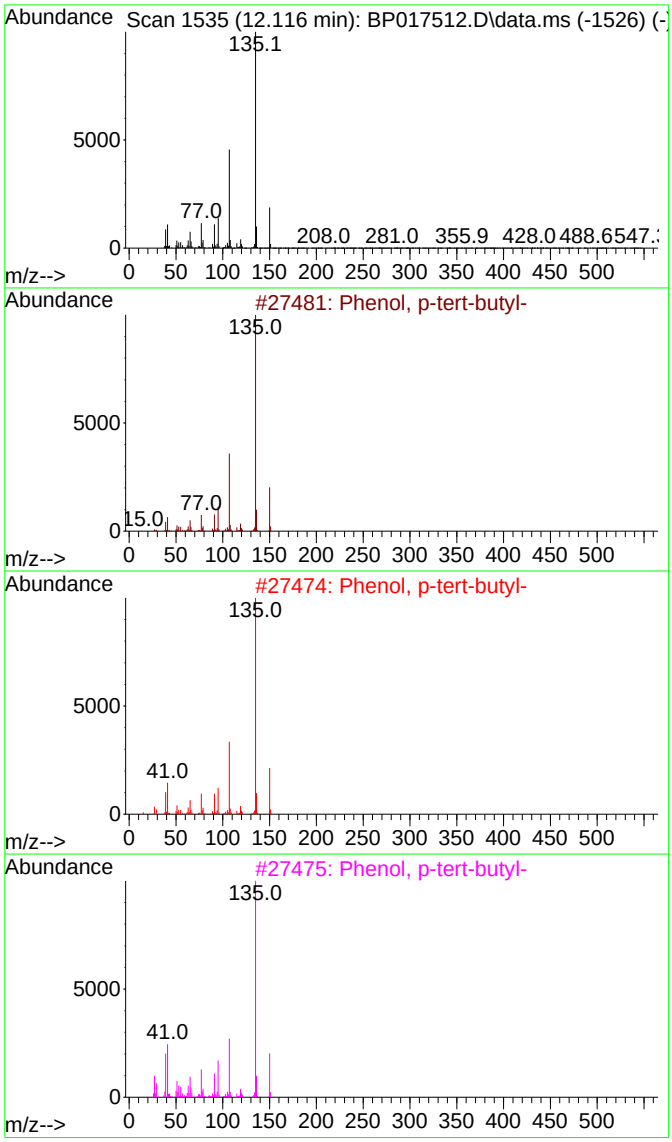
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

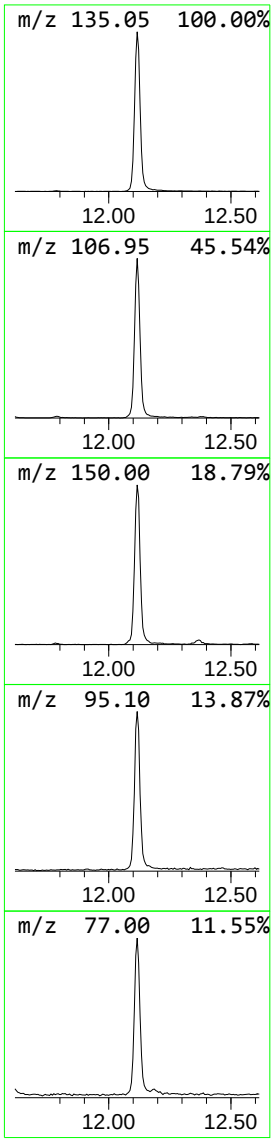
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.116	31.72 ng/ul	1493790	Naphthalene-d8	10.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	93
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

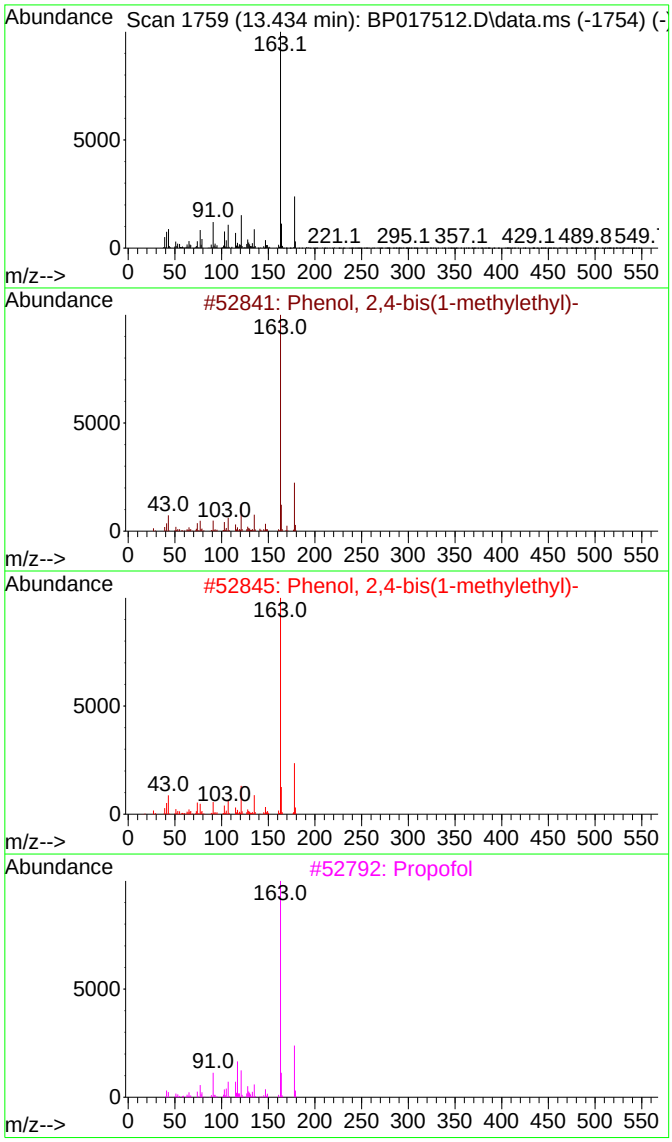
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

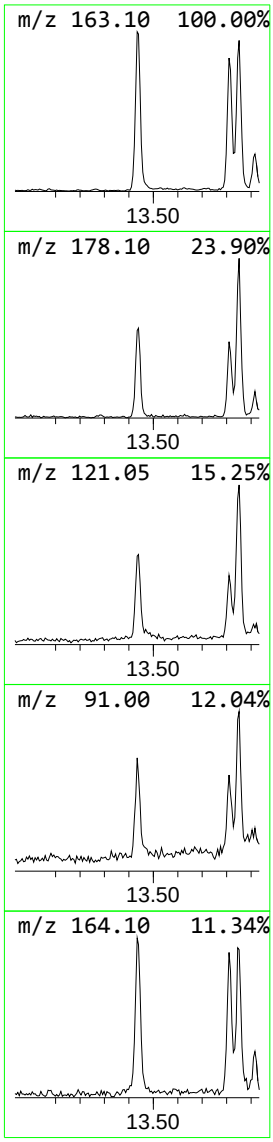
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Phenol, 2,4-bis(1-methyleth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.434	3.34 ng/ul	193936	Acenaphthene-d10	14.634

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	95
2	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	95
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	93
5	Propofol	178	C12H18O	002078-54-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

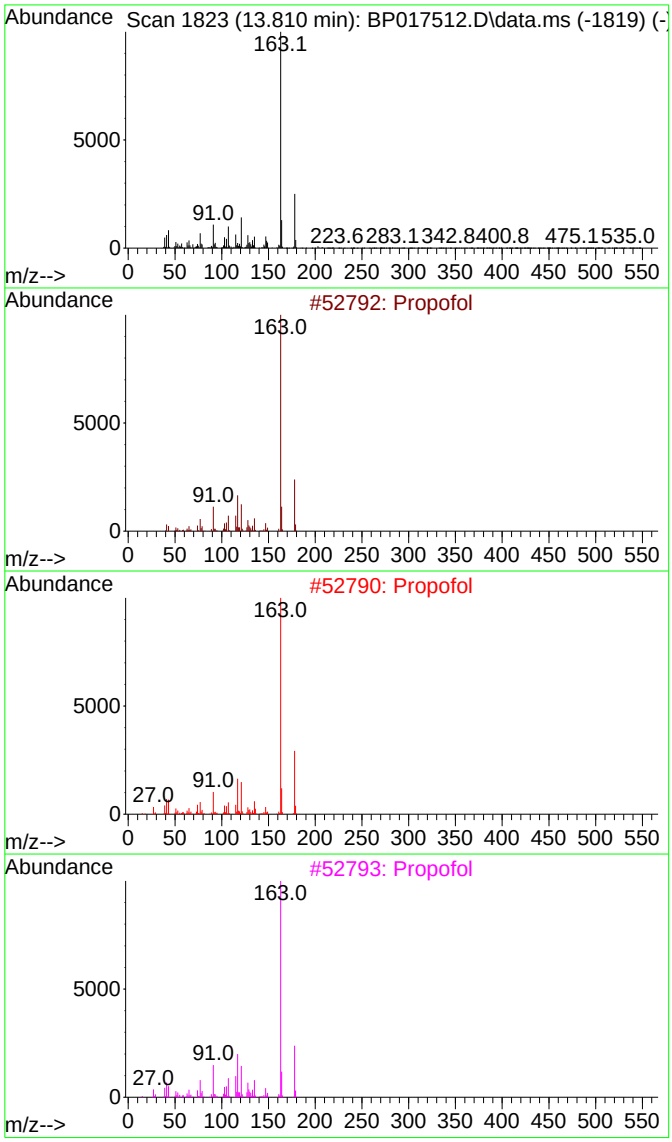
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

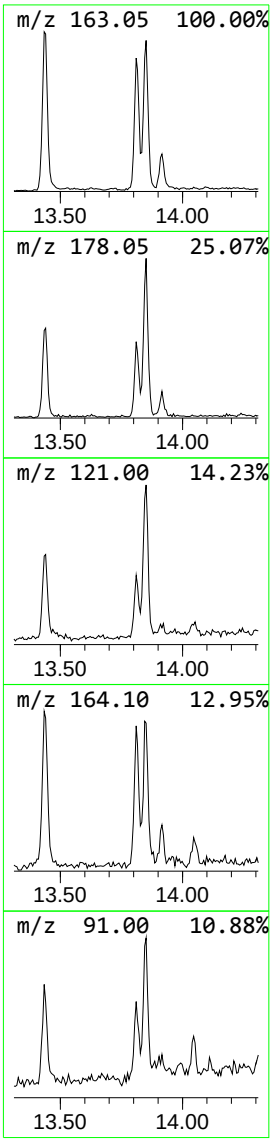
****

Peak Number 8 Propofol Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	2.39 ng/ul	138715	Acenaphthene-d10	14.634

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	94
2	Propofol	178	C12H18O	002078-54-8	94
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	91
5	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90





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Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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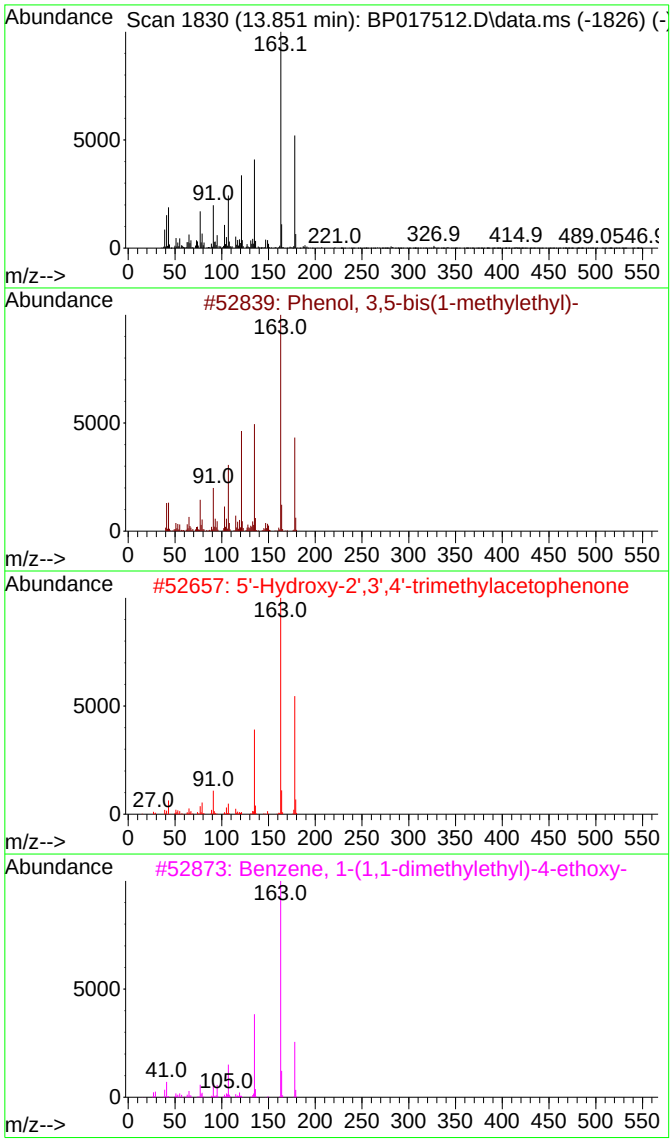
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

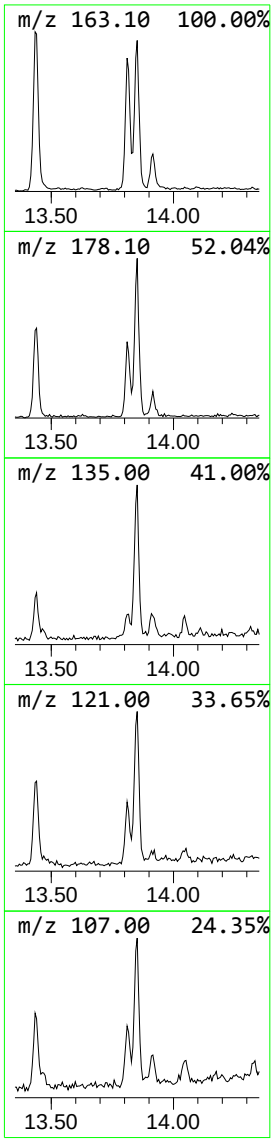
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 3,5-bis(1-methyleth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.851	4.12 ng/ul	239478	Acenaphthene-d10	14.634

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-bis(1-methylethyl)-	178	C12H18O	026886-05-5	93
2	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	87
3	Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	83
4	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	74
5	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	64





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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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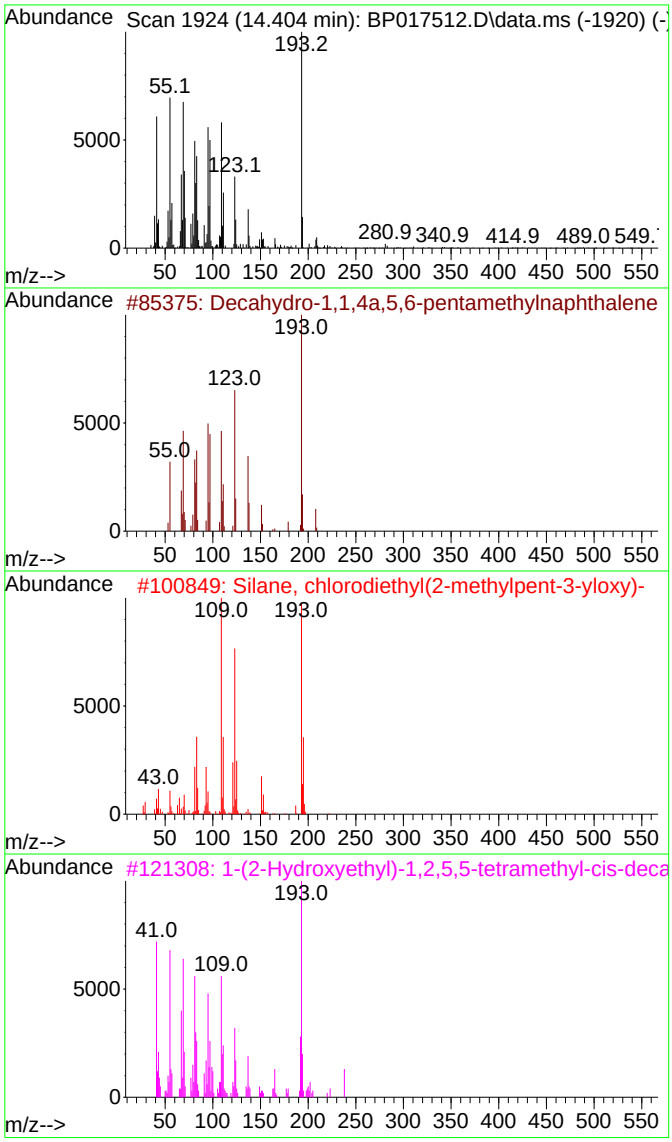
Quant Title : SVOA CALIBRATION

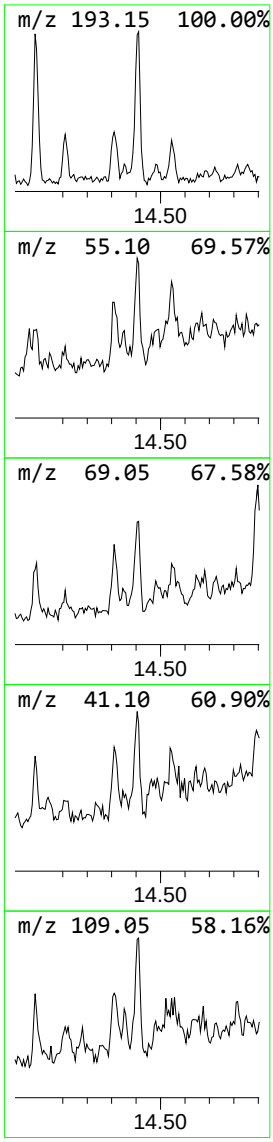
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.404	2.56 ng/ul	148864	Acenaphthene-d10	14.634

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	72
2	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	50
3	1-(2-Hydroxyethyl)-1,2,5,5-tetra...	238	C16H30O	1010298-97-4	42
4	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	35
5	3-((2R,5S)-7-(Cyclohexylsulfonyl...	340	C17H28N2O3S	1000483-82-6	35





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

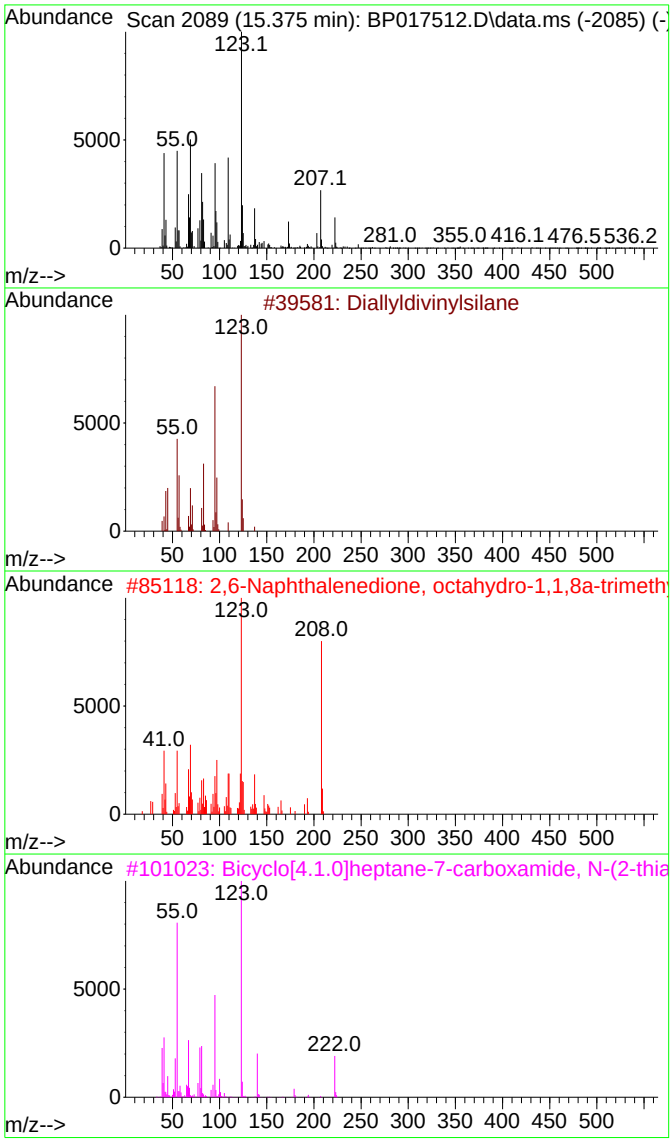
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

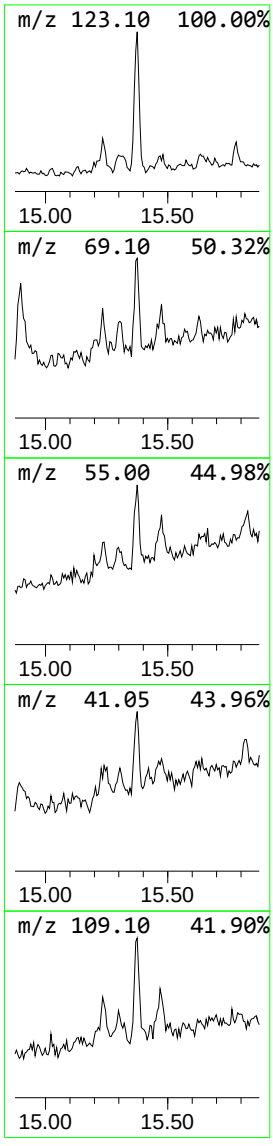
****

Peak Number 11 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.375	3.15 ng/ul	182793	Acenaphthene-d10	14.634

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diallyldivinylsilane	164	C10H16Si	119901-88-1	47
2	2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-16-4	47
3	Bicyclo[4.1.0]heptane-7-carboxam...	222	C11H14N2OS	329912-72-3	46
4	4-Nitrophenyl bicyclo[4.1.0]hept...	261	C14H15NO4	328938-65-4	43
5	2-Fluorobenzoic acid, 3-tetradec...	336	C21H33FO2	1000280-61-3	38





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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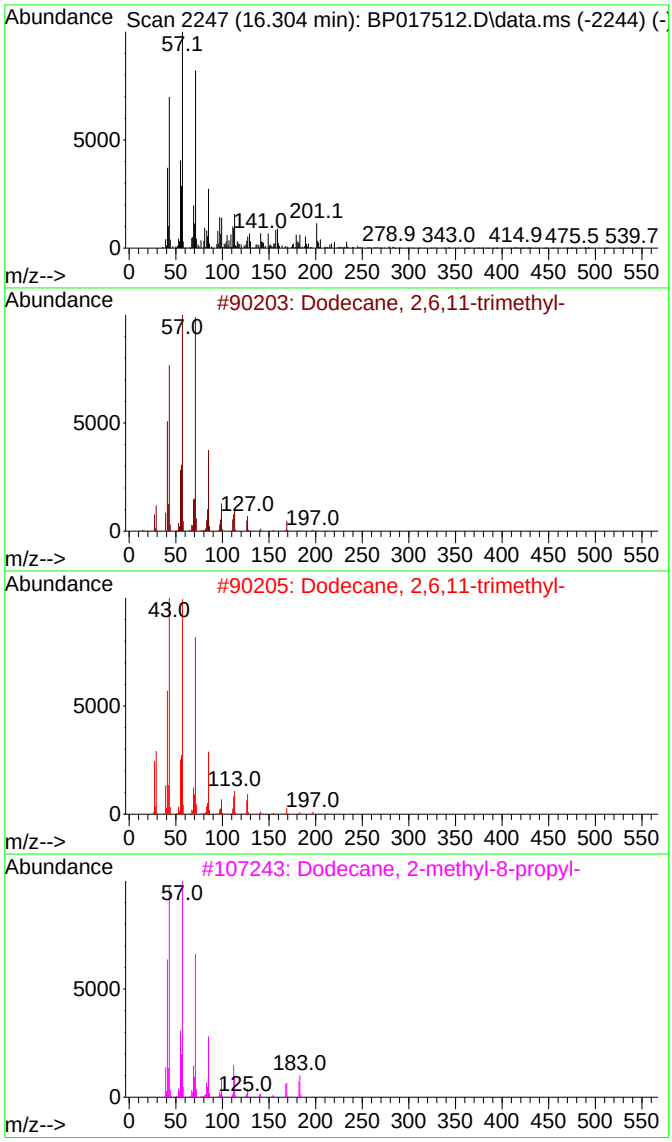
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

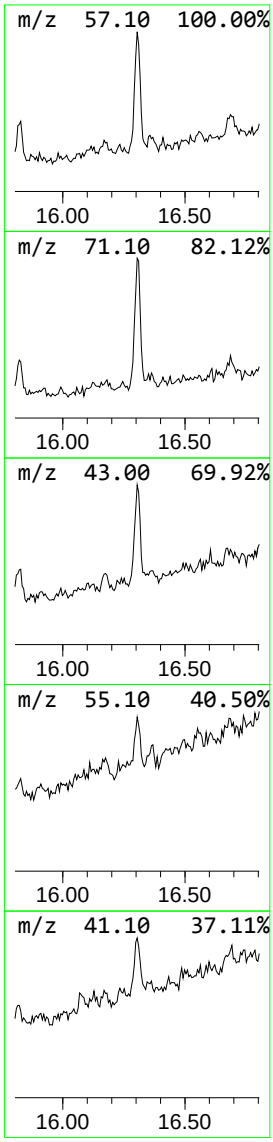
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.304	2.36 ng/ul	158992	Phenanthrene-d10	17.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	81
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80
3	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	74
4	Decane, 2-methyl-	156	C11H24	006975-98-0	74
5	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	72





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

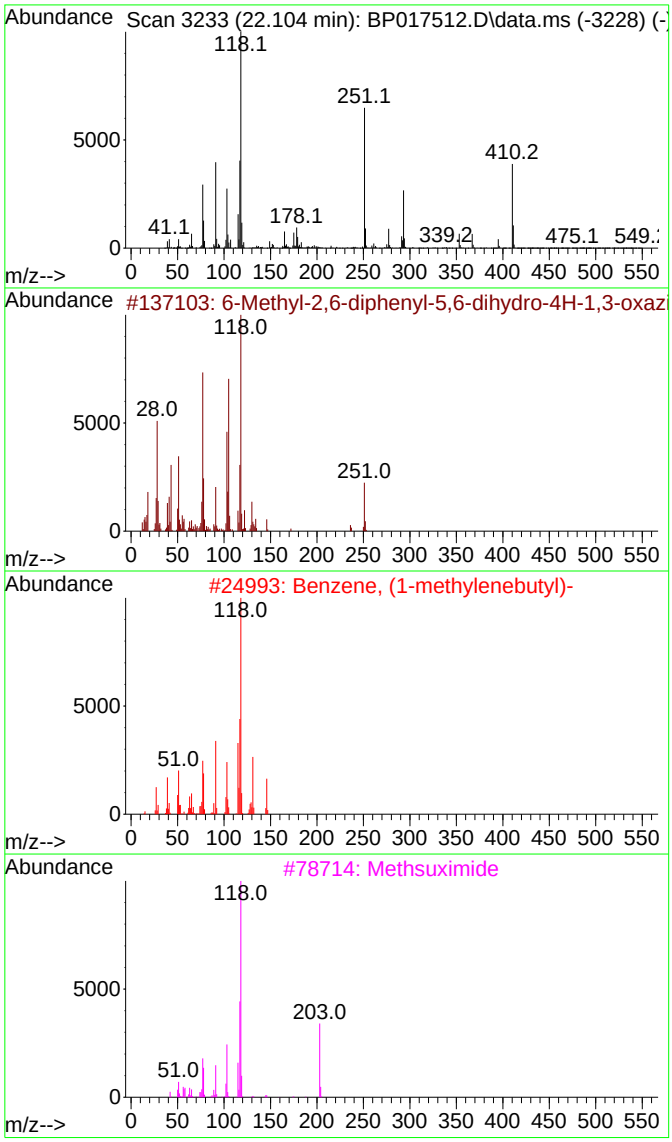
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

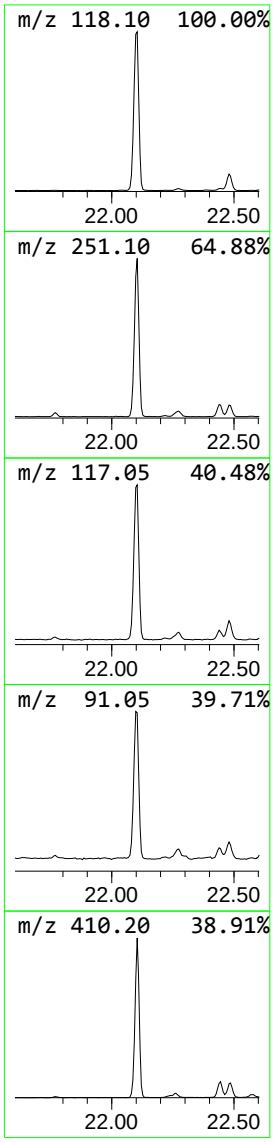
****

Peak Number 13 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.104	20.00 ng/ul	12038200	Chrysene-d12	21.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-2,6-diphenyl-5,6-dihydr...	251	C17H17NO	122035-87-4	49
2	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	38
3	Methsuximide	203	C12H13NO2	000077-41-8	38
4	Nicotinic acid, 3-phenylpropyl e...	241	C15H15NO2	1000308-37-0	38
5	Methsuximide	203	C12H13NO2	000077-41-8	35





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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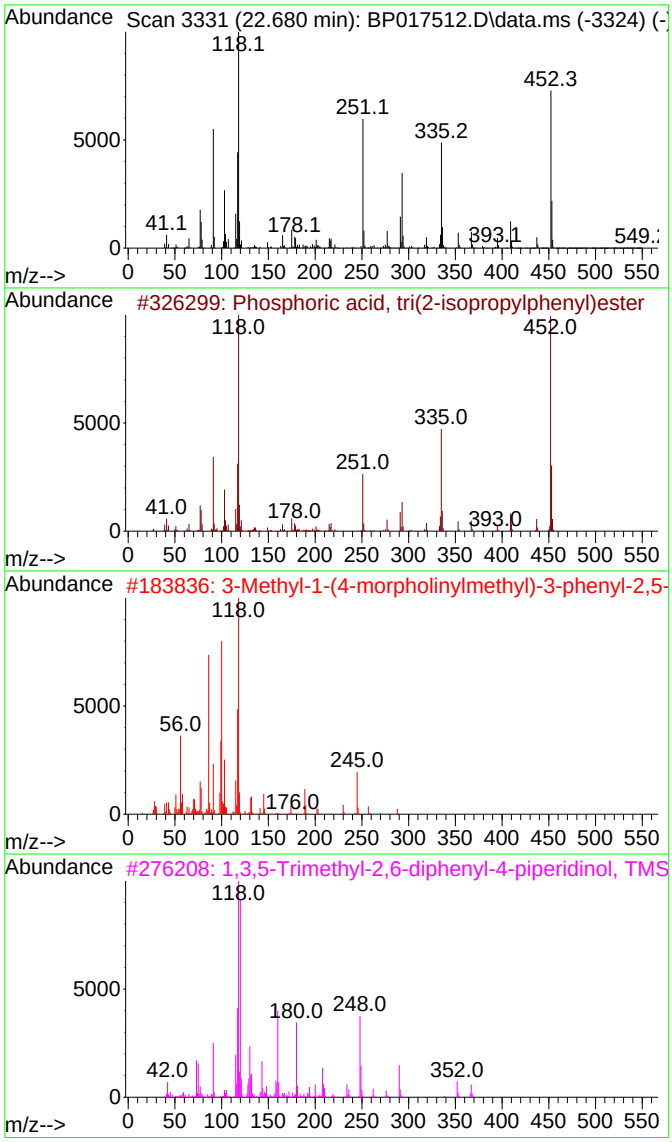
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

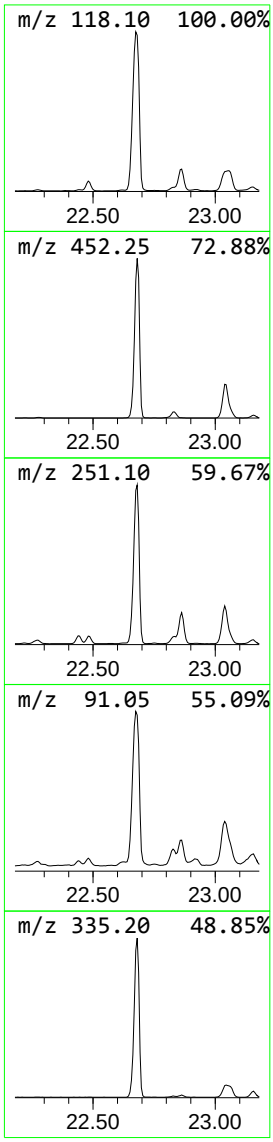
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Phosphoric acid, tri(2-isop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.680	50.96 ng/ul	30671600	Chrysene-d12	21.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, tri(2-isopropyl...	452	C27H33O4P	064532-95-2	98
2	3-Methyl-1-(4-morpholinylmethyl)...	288	C16H20N2O3	003780-72-1	38
3	1,3,5-Trimethyl-2,6-diphenyl-4-p...	367	C23H33NOSi	1000477-05-8	30
4	Bromoacetic acid, 3-phenylpropyl...	256	C11H13BrO2	059956-66-0	27
5	Diglycolic acid, pentyl 3-phenyl...	322	C18H26O5	1000382-17-3	27





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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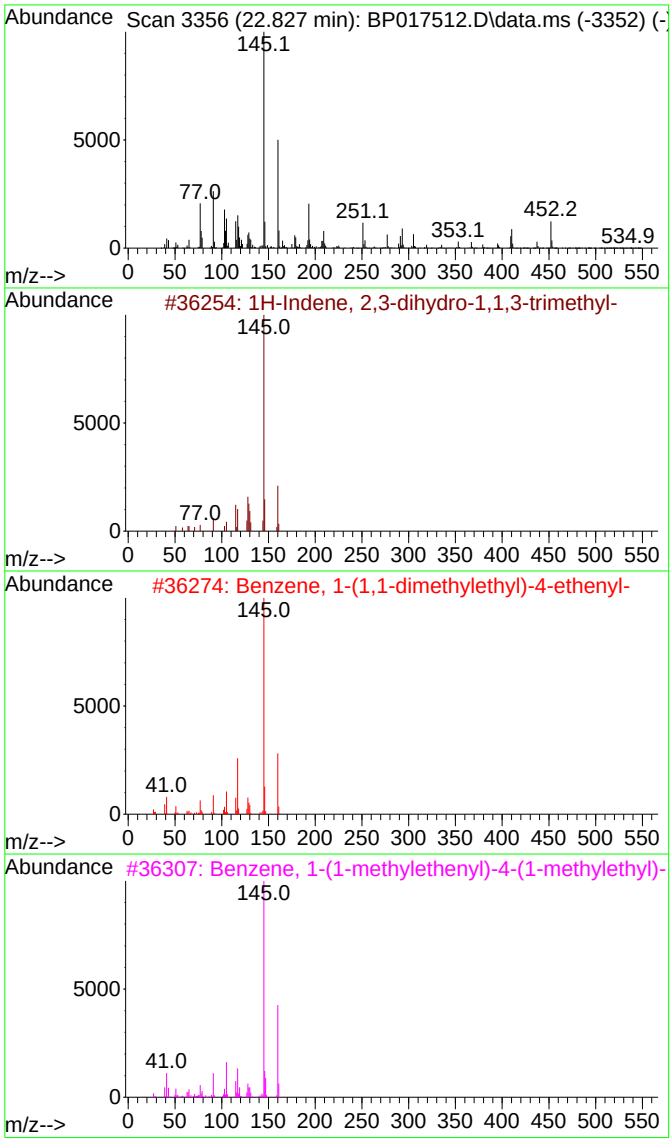
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

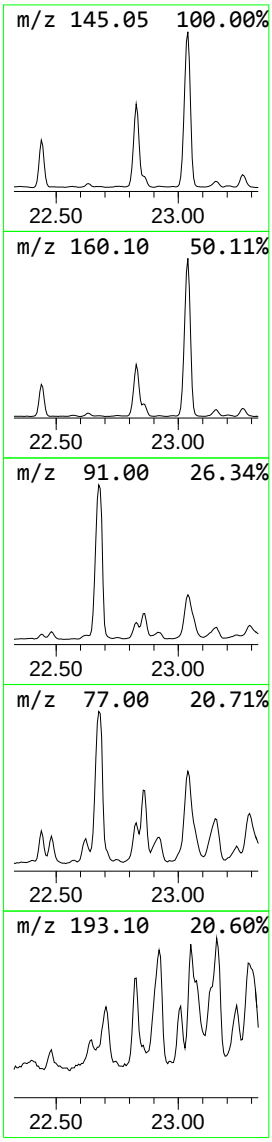
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.827	6.46 ng/ul	3887690	Chrysene-d12	21.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	60
2	Benzene, 1-(1,1-dimethylethyl)-4...	160	C12H16	001746-23-2	60
3	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	60
4	Benzene, 1-(1,1-dimethylethyl)-4...	160	C12H16	001746-23-2	60
5	Benzene, 1-(1,1-dimethylethyl)-4...	160	C12H16	001746-23-2	60





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Misc :
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

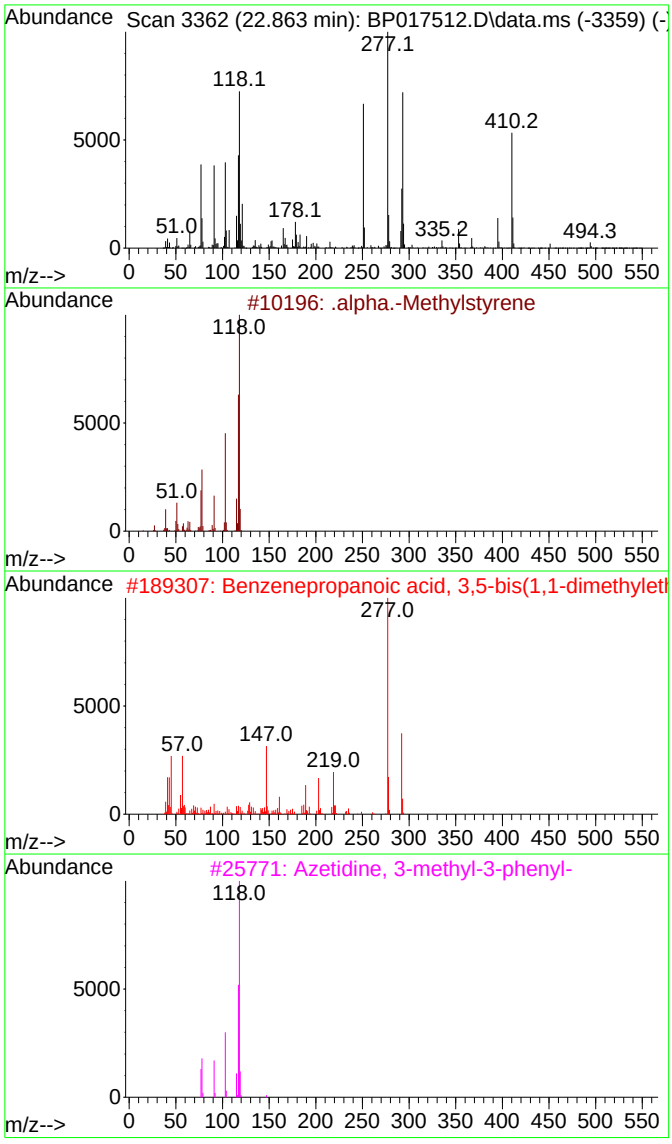
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

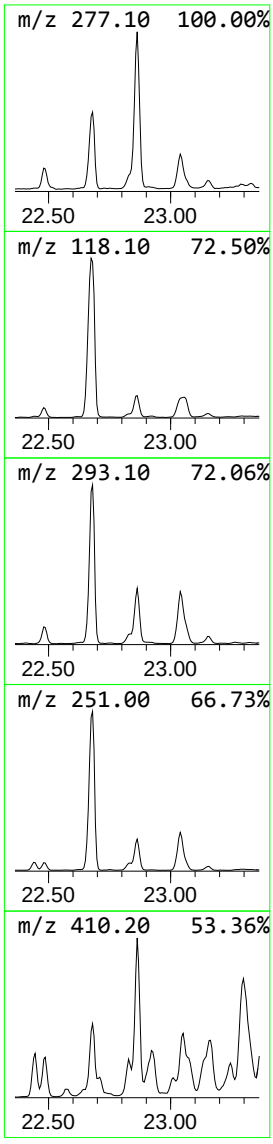
****

Peak Number 16 unknown-03 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.863	7.98 ng/ul	4803030	Chrysene-d12	21.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Methylstyrene	118	C9H10	000098-83-9	20
2	Benzenepropanoic acid, 3,5-bis(1...	292	C18H28O3	006386-38-5	14
3	Azetidine, 3-methyl-3-phenyl-	147	C10H13N	005961-33-1	14
4	.alpha.-Methylstyrene	118	C9H10	000098-83-9	14
5	Fumaric acid, 3-phenylpropyl 2,2...	348	C16H16F4O4	1000405-65-9	11





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Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

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ClientSampleId :
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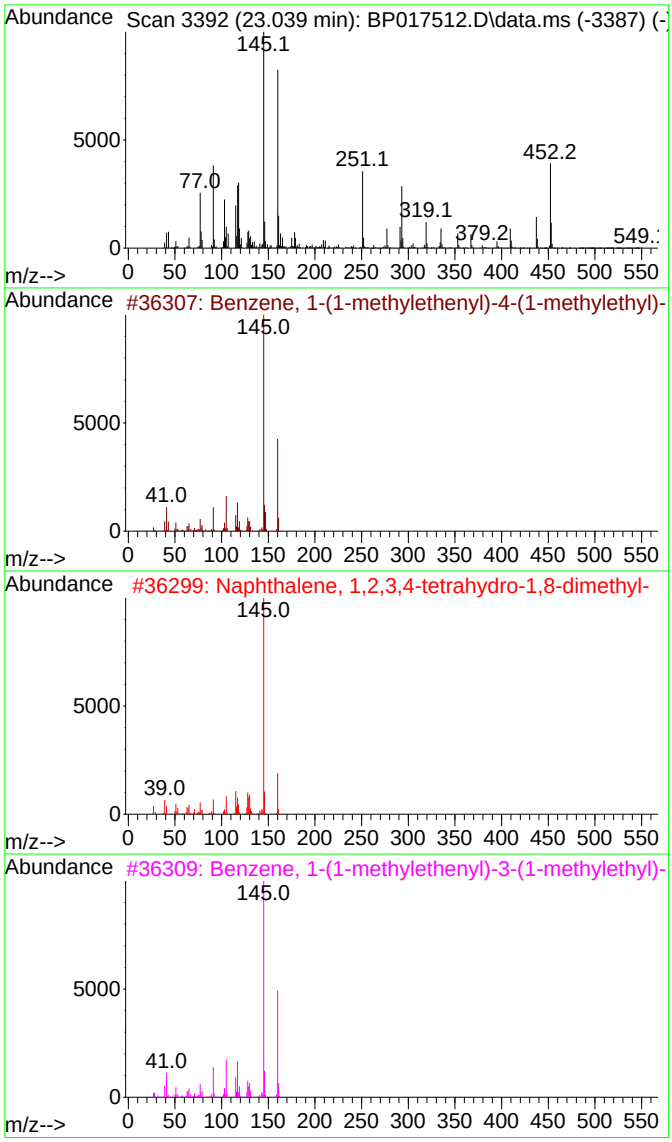
Quant Title : SVOA CALIBRATION

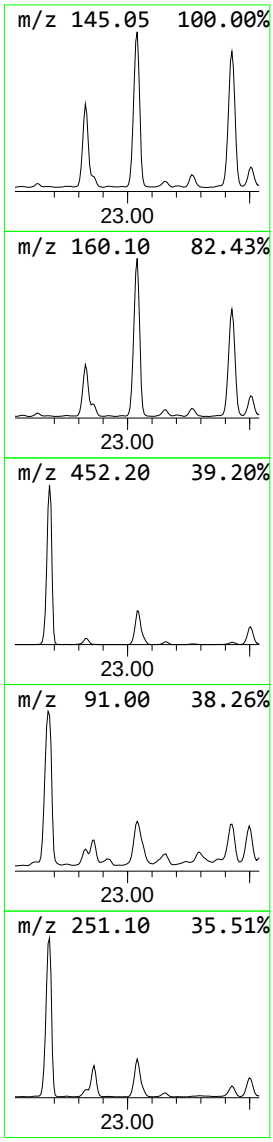
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-(1-methylethenyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.039	88.60 ng/ul	16196300	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	55
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	49
3	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	49
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	49
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	49





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Sample : 04417-30
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ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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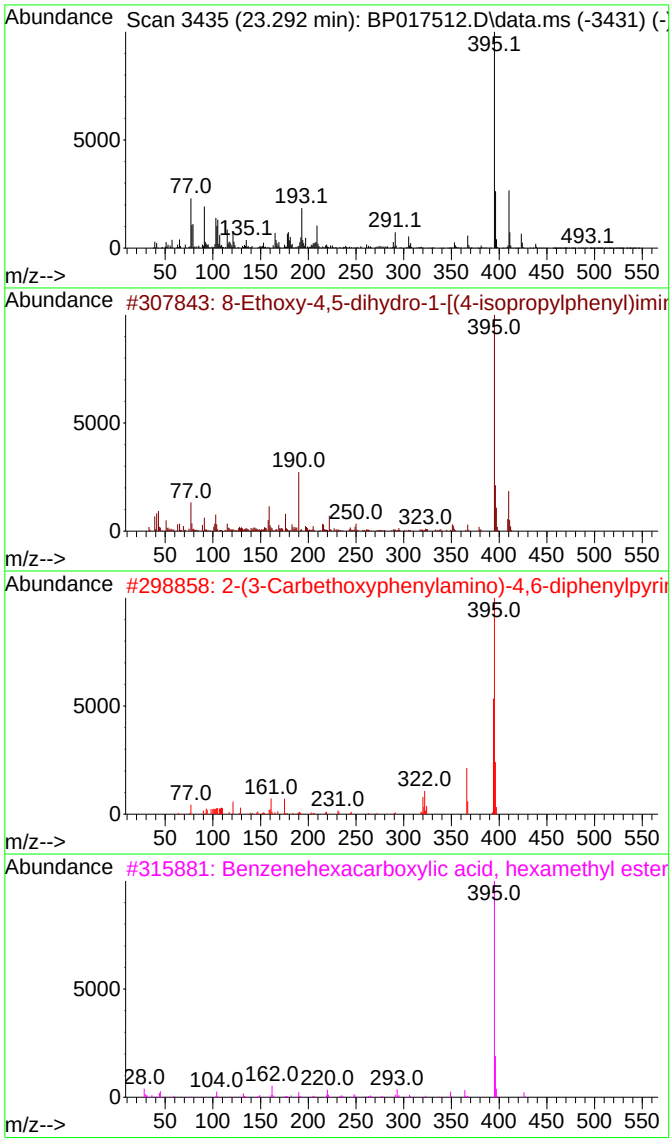
Quant Title : SVOA CALIBRATION

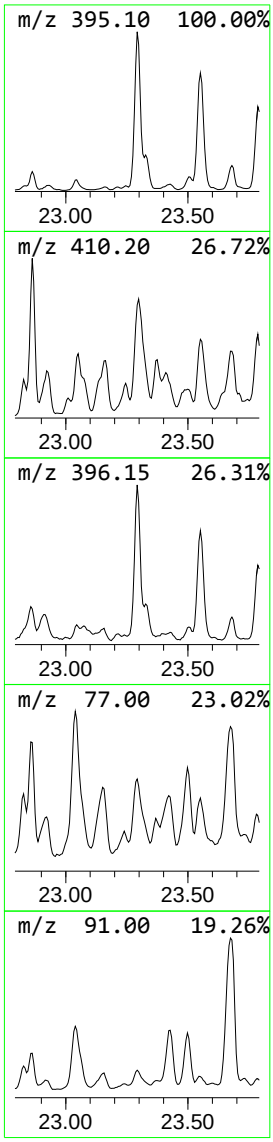
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 8-Ethoxy-4,5-dihydro-1-[(4-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.292	24.09 ng/ul	4404000	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Ethoxy-4,5-dihydro-1-[(4-isopr...	410	C23H26N2OS2	326174-54-3	72
2	2-(3-Carbethoxyphenylamino)-4,6-...	395	C25H21N3O2	1010070-62-1	58
3	Benzenehexacarboxylic acid, hexa...	426	C18H18O12	006237-59-8	53
4	N,N'-Bis(salicylidene)-3,3'-bis(...	395	C20H23N3NiO2	1010105-09-8	43
5	Phosphine imide, N-phenyl-P,P,P-...	395	C27H26NP	014796-90-8	42





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Quant Title : SVOA CALIBRATION

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

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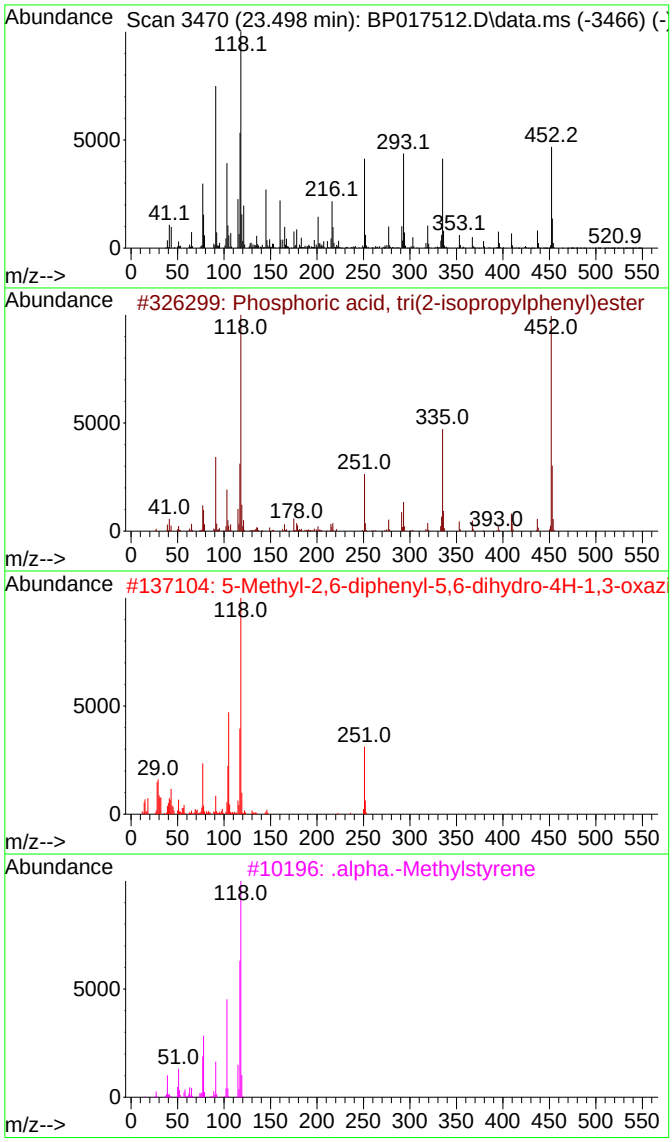
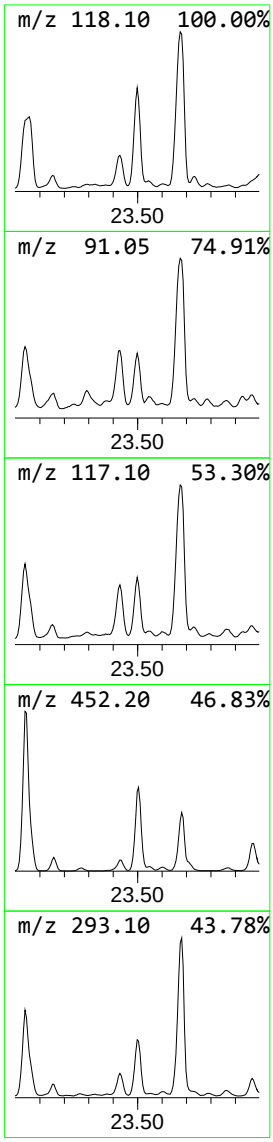
Peak Number 19 unknown-04 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.498	34.24 ng/ul	6259270	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, tri(2-isopropyl...	452	C27H33O4P	064532-95-2	91
2	5-Methyl-2,6-diphenyl-5,6-dihydr...	251	C17H17NO	023665-17-0	41
3	.alpha.-Methylstyrene	118	C9H10	000098-83-9	38
4	Nicotinic acid, 3-phenylpropyl e...	241	C15H15NO2	1000308-37-0	25
5	Isonicotinic acid, 3-phenylpropy...	241	C15H15NO2	1000308-36-4	22





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Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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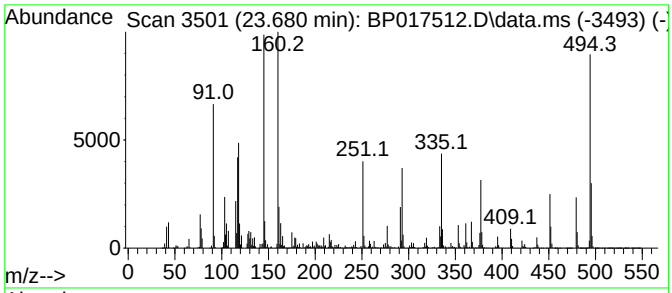
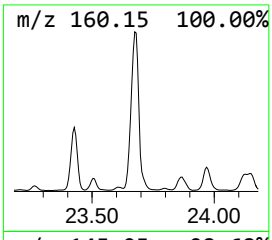
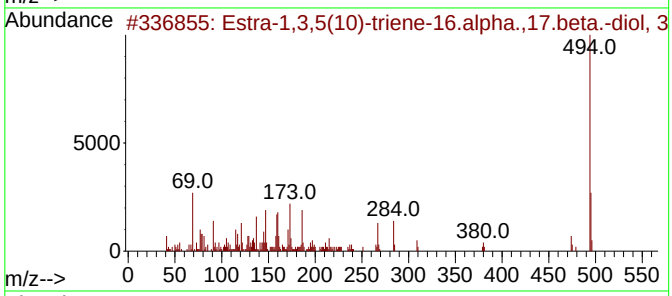
Peak Number 20 unknown-05 Concentration Rank 3

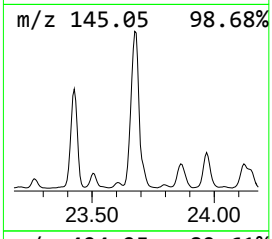
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.680	173.01 ng/ul	31625300	Perylene-d12	24.168

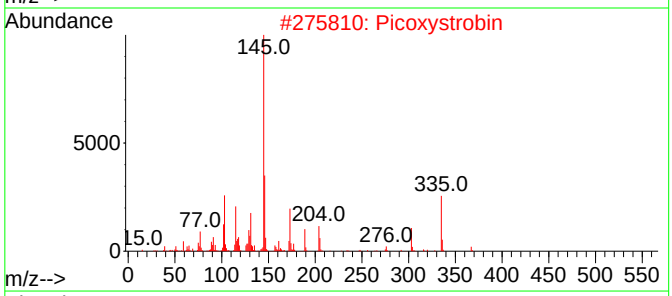
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Estra-1,3,5(10)-triene-16.alpha....	494	C23H24F6O5	034210-15-6	9
2	Picoxystrobin	367	C18H16F3NO4	117428-22-5	9
3	Cholestan-3-one, 2-(phenylthio)-...	494	C33H50OS	063608-48-0	9
4	1,4-Bis-(1,3,7-trimethyl-2,6-dio...	494	C22H22N8O6	1000286-01-3	9
5	Picoxystrobin	367	C18H16F3NO4	117428-22-5	9

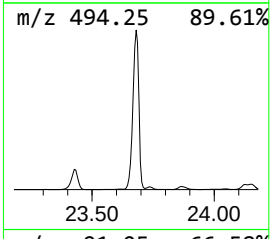


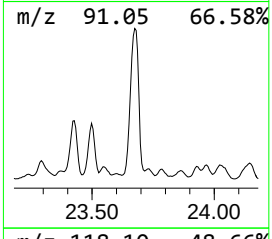


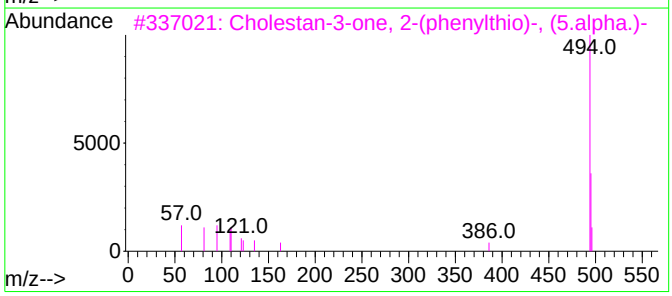


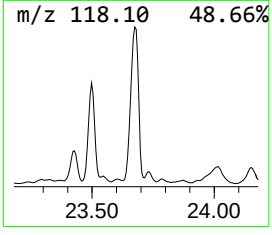












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Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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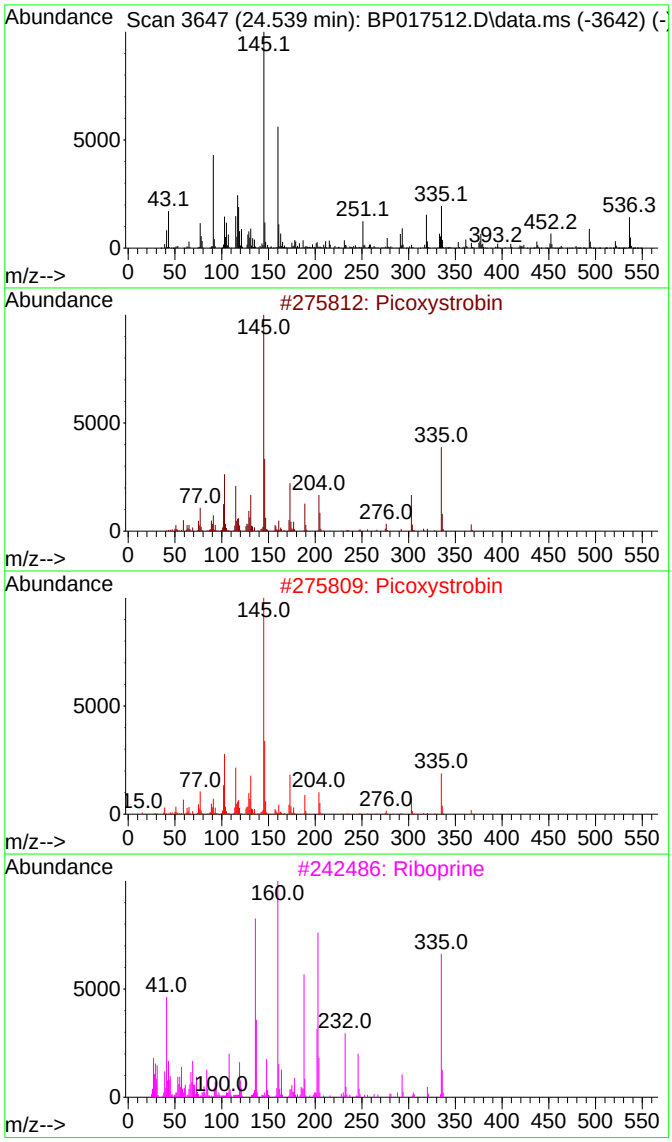
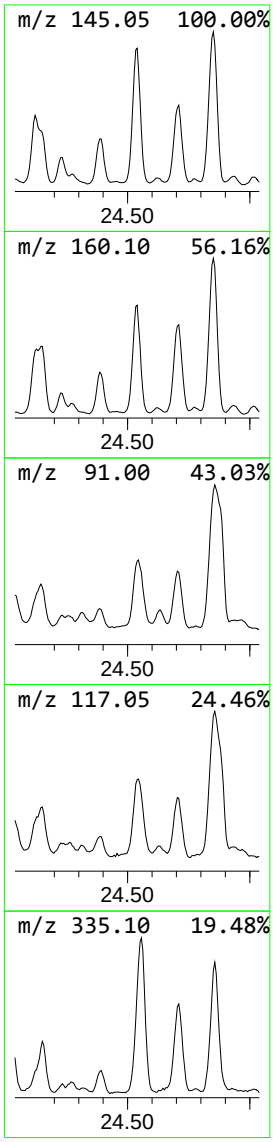
Peak Number 21 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	35.65 ng/ul	6515960	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Picoxystrobin	367	C18H16F3NO4	117428-22-5	25
2	Picoxystrobin	367	C18H16F3NO4	117428-22-5	25
3	Riboprine	335	C15H21N5O4	007724-76-7	22
4	Picoxystrobin	367	C18H16F3NO4	117428-22-5	10
5	Picoxystrobin	367	C18H16F3NO4	117428-22-5	10





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Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

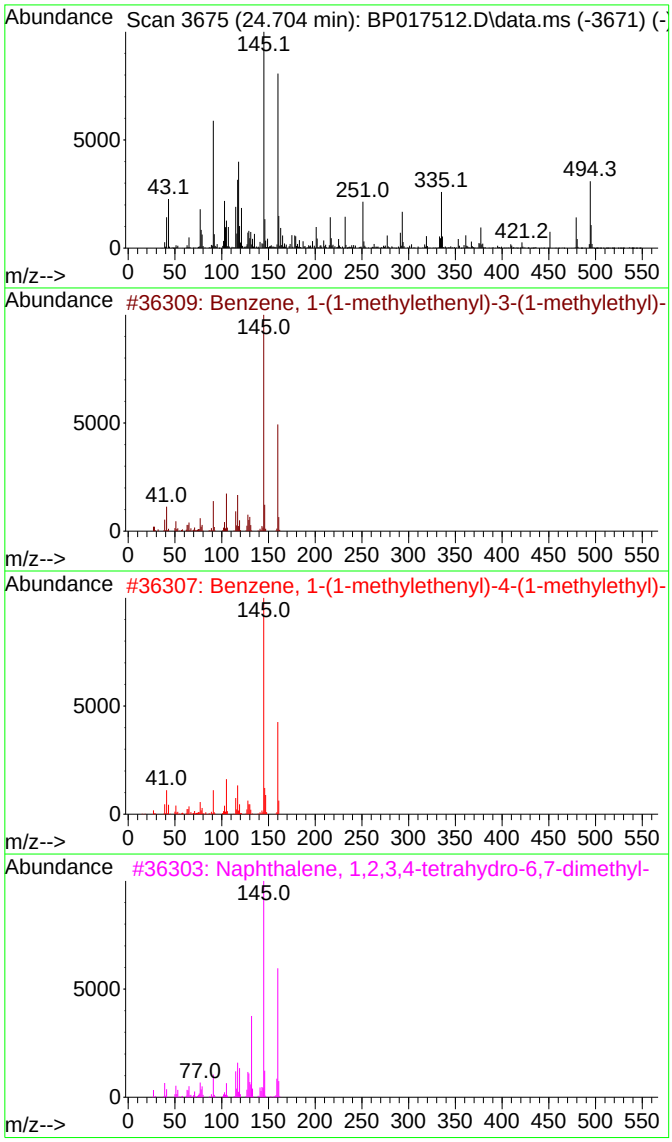
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

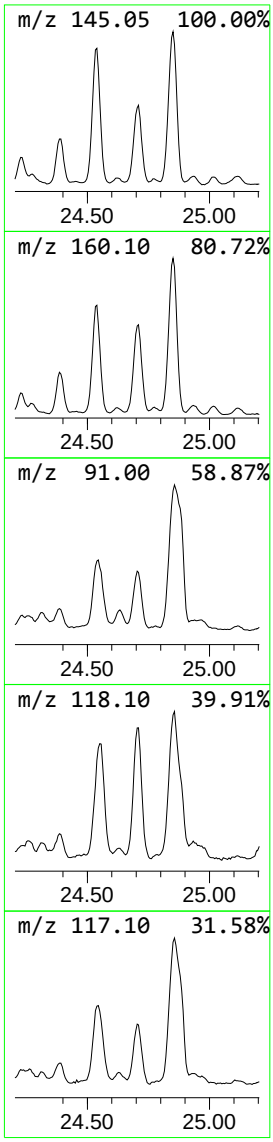
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Benzene, 1-(1-methylethenyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.704	28.36 ng/ul	5184620	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	55
2	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	55
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	49
4	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	49
5	1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	49





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Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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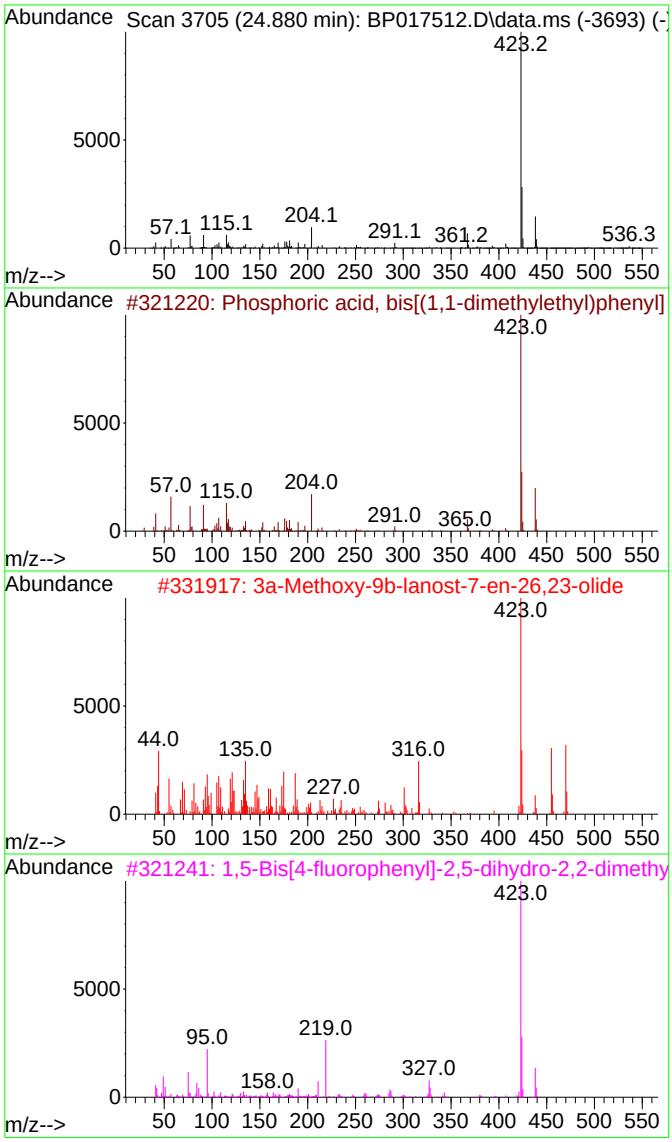
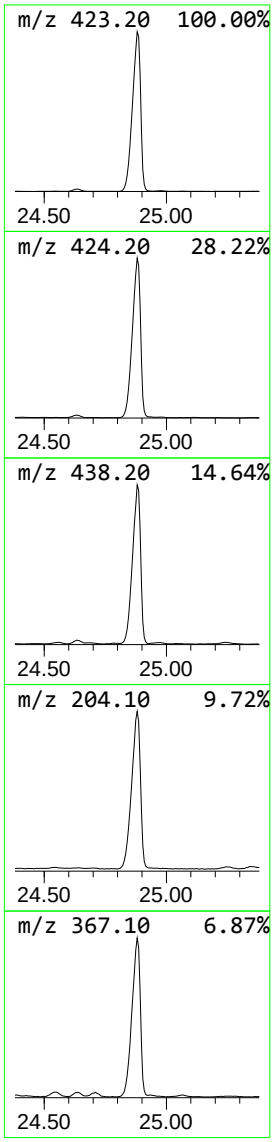
Peak Number 23 3a-Methoxy-9b-lanost-7-en-2... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.880	215.28 ng/ul	39351800	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, bis[(1,1-dimeth...	438	C26H31O4P	065652-41-7	98
2	3a-Methoxy-9b-lanost-7-en-26,23-...	470	C31H50O3	123073-95-0	72
3	1,5-Bis[4-fluorophenyl]-2,5-dihy...	438	C27H20F2N4	004447-85-2	56
4	Phenol, 2,3,5,6-tetrabromo-4-ethyl-	434	C8H6Br4O	105590-24-7	50
5	2-(4-Phenylpiperazino)-3-(4-toli...	423	C27H25N3O2	1000110-91-2	45





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Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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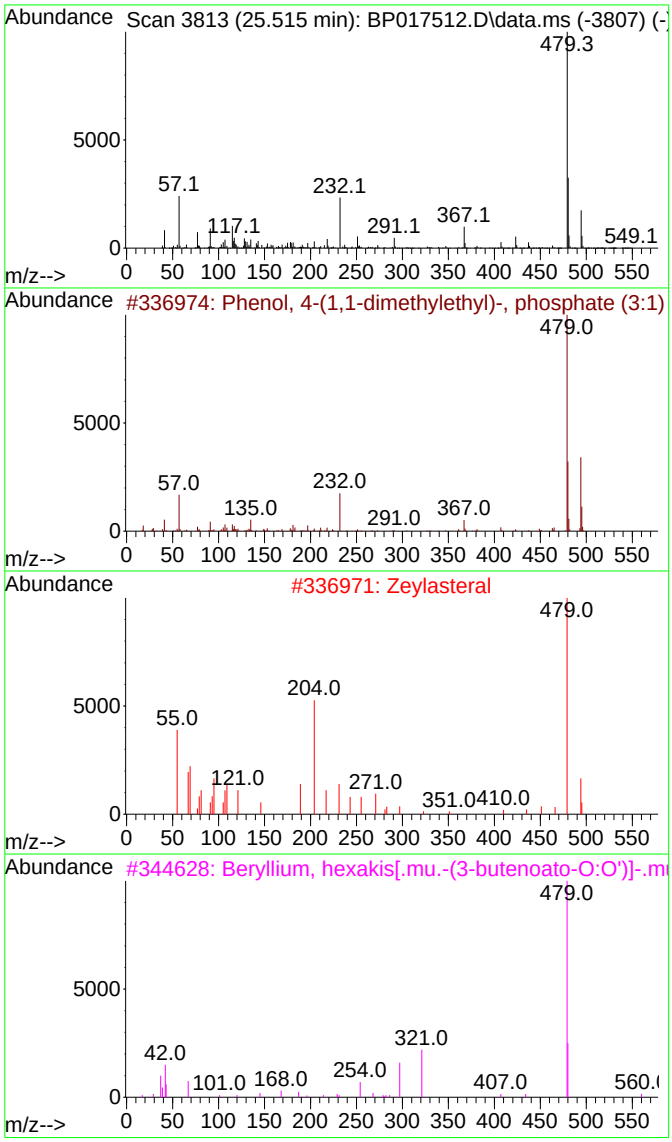
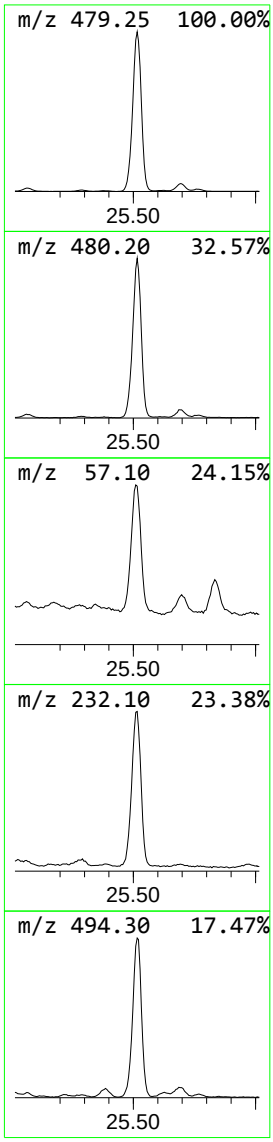
Peak Number 24 unknown-07 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.515	47.73 ng/ul	8724210	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylethyl)-, ...	494	C30H39O4P	000078-33-1	35
2	Zeylasteral	494	C30H38O6	087064-16-2	28
3	Beryllium, hexakis[.mu.-(3-buten...	562	C24H30Be4O13	056377-88-9	9
4	Dihydromorphine, O,O'-bis(triflu...	479	C21H19F6NO5	1000448-10-6	5
5	Silane, diethyl(dodec-9-ynyloxy)...	508	C32H64O2Si	1000363-58-8	4





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Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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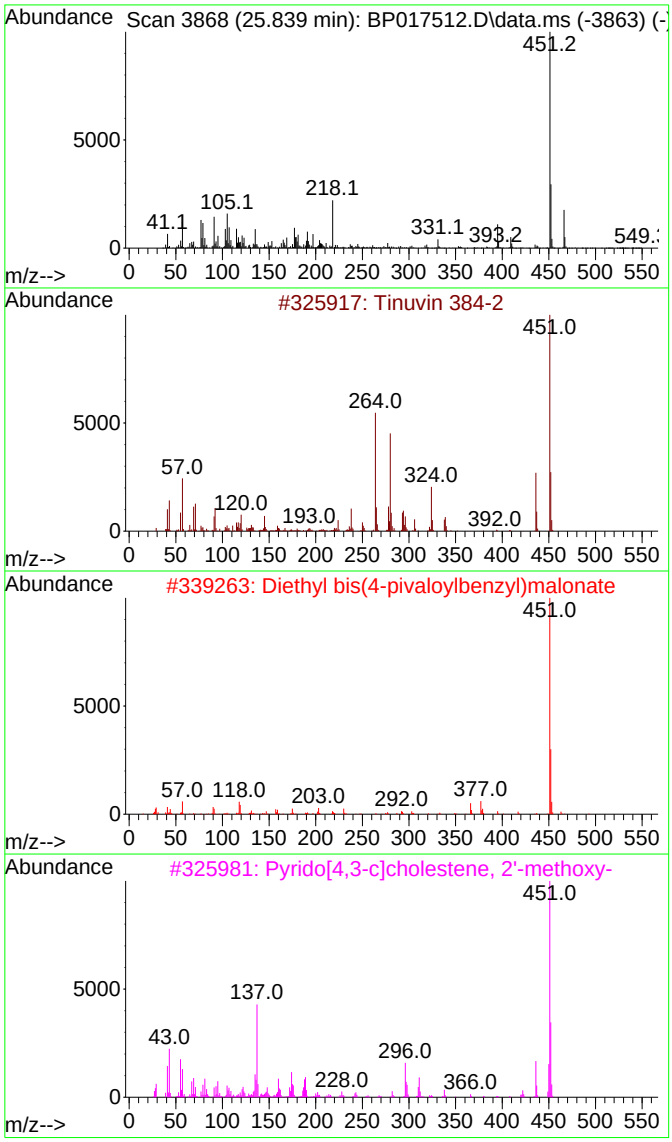
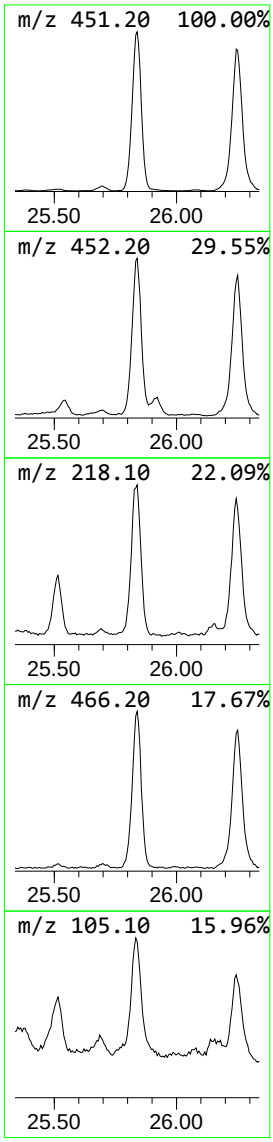
Peak Number 25 Tinuvin 384-2 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.839	32.95 ng/ul	6022700	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tinuvin 384-2	451	C27H37N3O3	084268-22-4	50
2	Diethyl bis(4-pivaloylbenzyl)mal...	508	C31H40O6	053780-45-3	40
3	Pyrido[4,3-c]cholestene, 2'-meth...	451	C31H49NO	1000210-91-7	40
4	Pregn-5-en-18-oic acid, 3-[(5,5-...	451	C29H41NO3	055401-43-9	38
5	Licochalcone C, 2TMS (isomer 2)	482	C27H38O4Si2	1000495-08-5	33





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Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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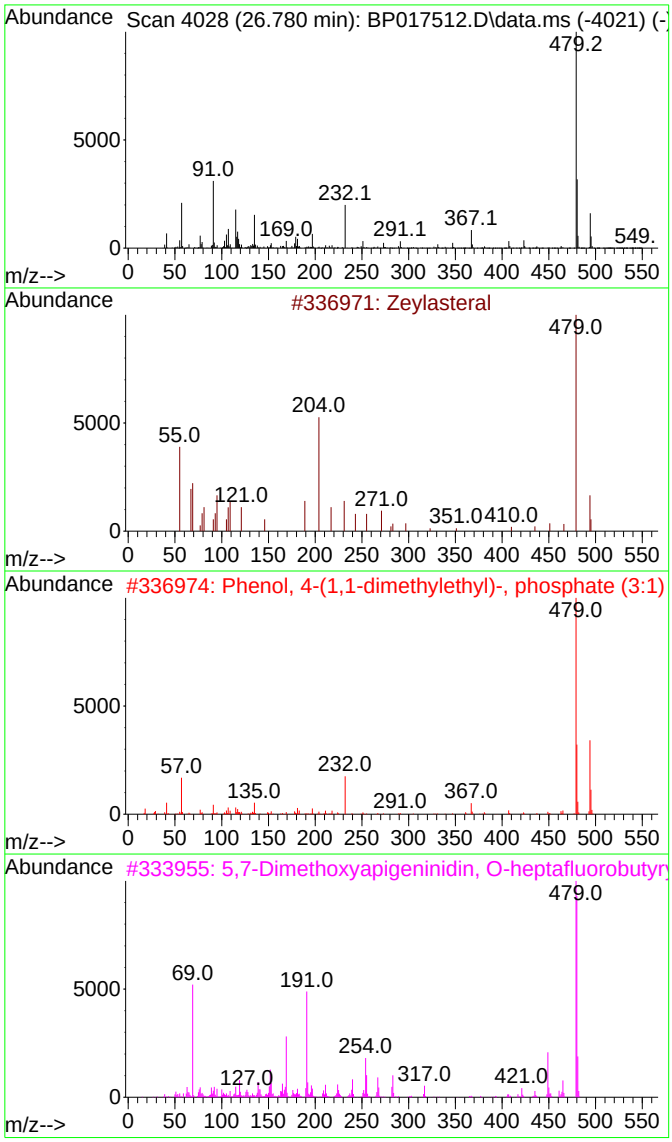
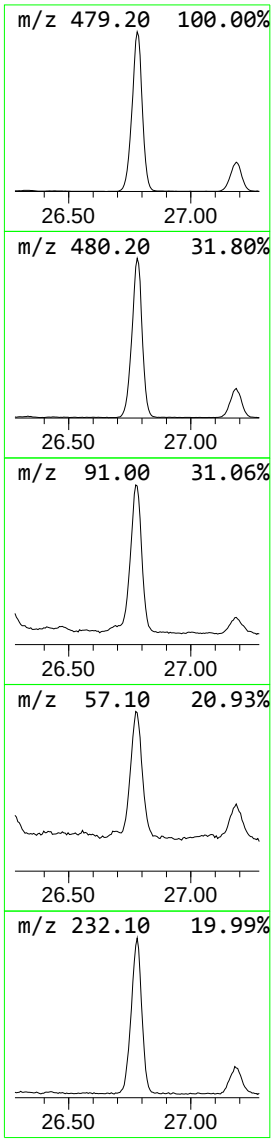
Peak Number 26 unknown-08 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.780	48.92 ng/ul	8943080	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Zeylasteral	494	C30H38O6	087064-16-2	37
2	Phenol, 4-(1,1-dimethylethyl)-, ...	494	C30H39O4P	000078-33-1	10
3	5,7-Dimethoxyapigeninidin, O-hep...	480	C21H15F7O5	1000453-04-1	4
4	Silane, methylvinyl(pentafluorob...	494	C25H39F5O2Si	1010420-62-1	2
5	Dihydromorphine, O,O'-bis(triflu...	479	C21H19F6NO5	1000448-10-6	1





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Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

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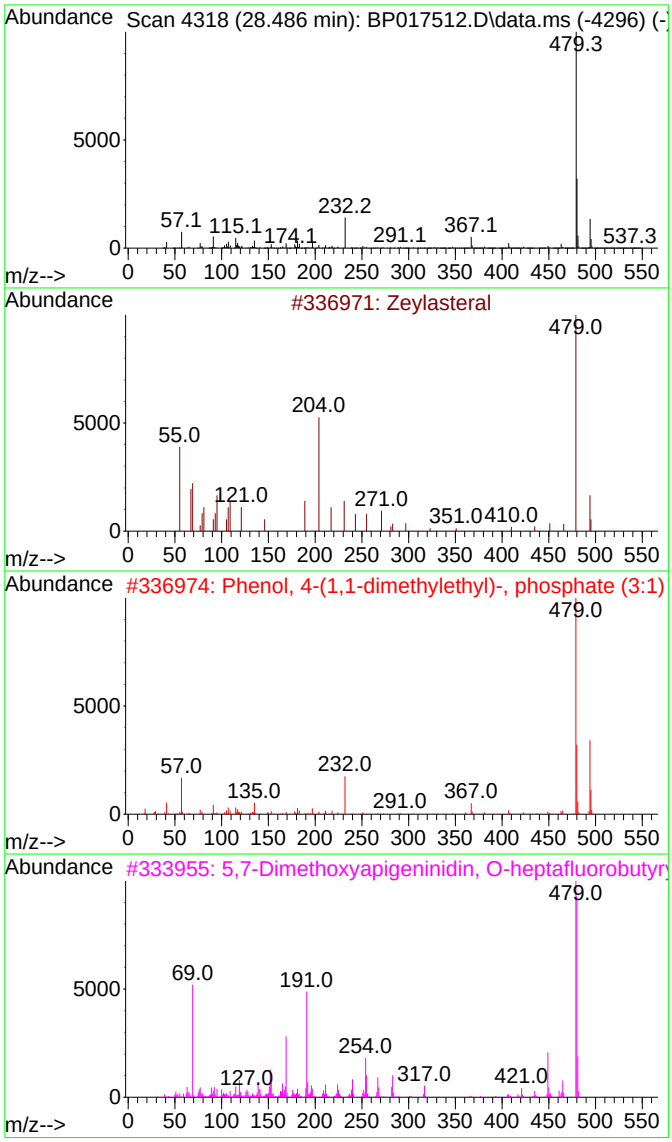
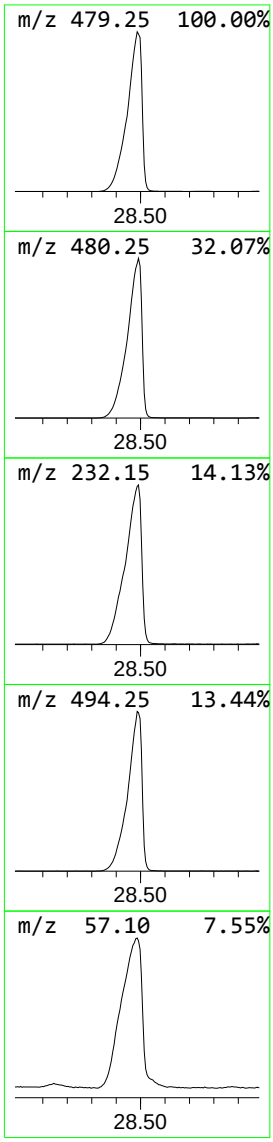
Peak Number 27 unknown-09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.486	371.50 ng/ul	67908600	Perylene-d12	24.168

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Zeylasteral	494	C30H38O6	087064-16-2	33
2	Phenol, 4-(1,1-dimethylethyl)-, ...	494	C30H39O4P	000078-33-1	30
3	5,7-Dimethoxyapigeninidin, O-hep...	480	C21H15F7O5	1000453-04-1	9
4	Beryllium, hexakis[.mu.-(3-buten...	562	C24H30Be4O13	056377-88-9	9
5	Silane, diethyl(dodec-9-ynyloxy)...	508	C32H64O2Si	1000363-58-8	2





Tentatively Identified Compound (LSC) summary

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Data File : BP017512.D
Acq On : 03 Oct 2023 12:07
Operator : MA/JU
Sample : 04417-30
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYM9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.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
Tetrahydropyran...	3.834	3.3	ng/ul	117238	1	7.934	720771	20.0
Butane, 2,3-dim...	3.911	2.8	ng/ul	100288	1	7.934	720771	20.0
Phenol, 2,5-dim...	10.216	2.1	ng/ul	100384	2	10.775	941929	20.0
Phenol, 2-(1-me...	10.681	4.0	ng/ul	189683	2	10.775	941929	20.0
p-Cumenol	11.122	2.7	ng/ul	126445	2	10.775	941929	20.0
Phenol, p-tert-...	12.116	31.7	ng/ul	1493790	2	10.775	941929	20.0
Phenol, 2,4-bis...	13.434	3.3	ng/ul	193936	3	14.634	1162310	20.0
Propofol	13.810	2.4	ng/ul	138715	3	14.634	1162310	20.0
Phenol, 3,5-bis...	13.851	4.1	ng/ul	239478	3	14.634	1162310	20.0
Decahydro-1,1,4...	14.404	2.6	ng/ul	148864	3	14.634	1162310	20.0
unknown-01	15.375	3.1	ng/ul	182793	3	14.634	1162310	20.0
(DEL) Alkane: S...	16.304	2.4	ng/ul	158992	4	17.422	1347940	20.0
unknown-02	22.104	20.0	ng/ul	12038200	5	21.569	12038200	20.0
Phosphoric acid...	22.680	51.0	ng/ul	30671600	5	21.569	12038200	20.0
1H-Indene, 2,3-...	22.827	6.5	ng/ul	3887690	5	21.569	12038200	20.0
unknown-03	22.863	8.0	ng/ul	4803030	5	21.569	12038200	20.0
Benzene, 1-(1-m...	23.039	88.6	ng/ul	16196300	6	24.168	3655870	20.0
8-Ethoxy-4,5-di...	23.292	24.1	ng/ul	4404000	6	24.168	3655870	20.0
unknown-04	23.498	34.2	ng/ul	6259270	6	24.168	3655870	20.0
unknown-05	23.680	173.0	ng/ul	31625300	6	24.168	3655870	20.0
unknown-06	24.539	35.6	ng/ul	6515960	6	24.168	3655870	20.0
Benzene, 1-(1-m...	24.704	28.4	ng/ul	5184620	6	24.168	3655870	20.0
3a-Methoxy-9b-l...	24.880	215.3	ng/ul	39351800	6	24.168	3655870	20.0
unknown-07	25.515	47.7	ng/ul	8724210	6	24.168	3655870	20.0
Tinuvin 384-2	25.839	33.0	ng/ul	6022700	6	24.168	3655870	20.0
unknown-08	26.780	48.9	ng/ul	8943080	6	24.168	3655870	20.0
unknown-09	28.486	371.5	ng/ul	67908600	6	24.168	3655870	20.0