LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

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-BM040922.M
Title : SVOA CALIBRATION

Signal : TIC: BM034681.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.025	672	678	696	rBV	368421	679796	0.19%	0.095%
2	7.190	700	706	718	rBV	377041	608057	0.17%	0.085%
3	7.390	734	740	750	rBV	450720	760923	0.21%	0.106%
4	7.860	814	820	831	rBV	508331	871479	0.24%	0.122%
5	8.578	933	942	955	rBV	344244	656017	0.18%	0.091%
6	9.025	1011	1018	1030	rBV	437488	767785	0.21%	0.107%
7	9.754	1135	1142	1156	rBV	321324	598802	0.16%	0.083%
8	10.296	1227	1234	1249	rBV	399445	801755	0.22%	0.112%
9	10.672	1290	1298	1313	rBV	681794	1181477	0.32%	0.165%
10	10.813	1315	1322	1344	rBV	279821	698410	0.19%	0.097%
11	13.166	1713	1722	1732	rBV	900215	1345890	0.37%	0.188%
12	13.878	1835	1843	1847	rBV	3195751	4464890	1.22%	0.623%
13	13.919	1847	1850	1855	rVV	1033316	1522816	0.42%	0.212%
14	14.201	1889	1898	1910	rBV	1162700	1754188	0.48%	0.245%
15	14.507	1944	1950	1959	rBV	948579	1451113	0.40%	0.202%
16	14.672	1971	1978	1984	rBV	3702128	5060026	1.38%	0.706%
17	15.501	2109	2119	2129	rVV	1488231	2216826	0.61%	0.309%
18	15.619	2135	2139	2145	rVV	232502	383765	0.10%	0.054%
19	15.683	2145	2150	2154	rVV	383209	515388	0.14%	0.072%
20	15.736	2154	2159	2169	rVB	595139	819040	0.22%	0.114%
21	16.172	2226	2233	2239	rBV	758325	1063007	0.29%	0.148%
22	16.283	2245	2252	2261	rVB	4109264	5267424	1.44%	0.734%
23	17.254	2411	2417	2422	rBV	1042645	1537456	0.42%	0.214%
24	17.354	2428	2434	2446	rVB	1272941	1972481	0.54%	0.275%
25	19.401	2775	2782	2787	rBV2	1082496	1664211	0.45%	0.232%
26	19.554	2803	2808	2811	rBV	1288213	1684587	0.46%	0.235%
27	19.583	2811	2813	2815	rVV	726788	913708	0.25%	0.127%
28	19.642	2819	2823	2827	rVV	1733617	2565448	0.70%	0.358%
29	19.677	2827	2829	2833	rVV3	577408	854976	0.23%	0.119%
30	19.724	2833	2837	2842	rVB2	1114228	1500858	0.41%	0.209%
31	19.789	2842	2848	2855	rBV	10060192	12153080	3.32%	1.695%
32	19.895	2860	2866	2870	rVV	8818278	11905480	3.25%	1.660%
33	19.936	2870	2873	2879	rVB	1916164	2294794	0.63%	0.320%
34	20.089	2893	2899	2901	rBV	1352698	2203913	0.60%	0.307%
35	20.124	2901	2905	2913	rVB	4844906	7203316	1.97%	1.004%
36	20.260	2924	2928	2932	rVB	471850	679339	0.19%	0.095%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

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-BM040922.M
Title : SVOA CALIBRATION

37	20.365	2933	2946	2955	rBV2	44864168	203658872	55.62%	28.397%
38	20.530	2968	2974	2988	rBV	12031477	27387029	7.48%	3.819%
39	20.760	3004	3013	3036	rBV	44384850	366156177	100.00%	51.055%
40	20.901	3036	3037	3039	rVB	35878864	12715342	3.47%	1.773%
41	21.160	3076	3081	3087	rVB2	430645	828059	0.23%	0.115%
42	21.307	3103	3106	3111	rVB2	348637	391892	0.11%	0.055%
43	21.454	3127	3131	3135	rBV	1125859	1347679	0.37%	0.188%
44	21.889	3201	3205	3210	rVV	537434	726789	0.20%	0.101%
45	21.954	3210	3216	3222	rVV	10678570	13276042	3.63%	1.851%
46	22.077	3231	3237	3242	rVB	1744671	2483166	0.68%	0.346%
47	22.424	3292	3296	3302	rVB2	248432	389771	0.11%	0.054%
48	22.607	3323	3327	3335	rVB	393370	624350	0.17%	0.087%
49	22.989	3388	3392	3398	rVB	237068	371474	0.10%	0.052%
50	23.130	3413	3416	3425	rVB2	243137	404852	0.11%	0.056%
51	23.654	3499	3505	3512	rBV	1028094	2031011	0.55%	0.283%
52	23.806	3525	3531	3553	rVB2	770494	1765753	0.48%	0.246%

Sum of corrected areas: 717180779

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Instrument :
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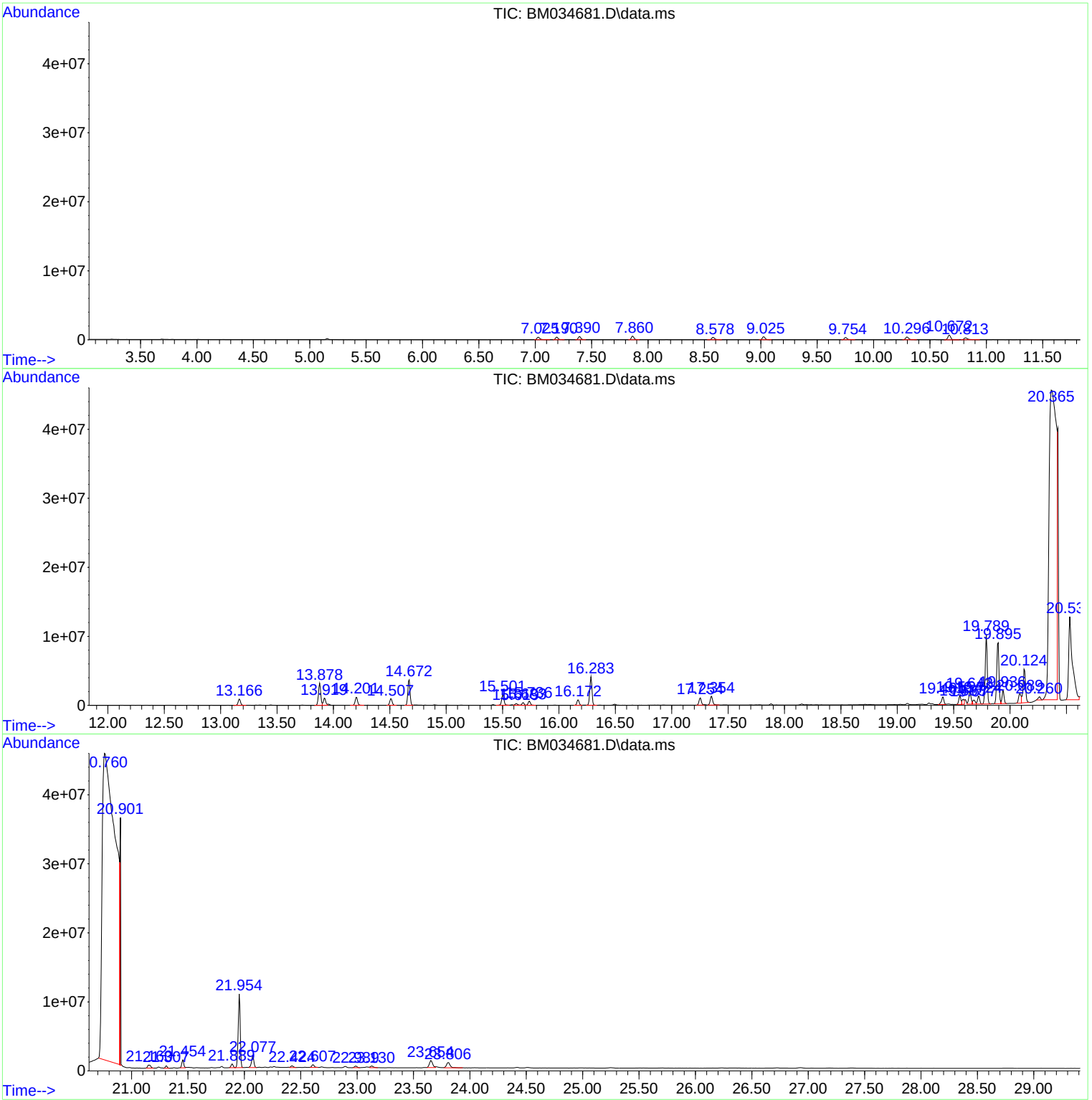
ClientSampleId :
ETNR3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
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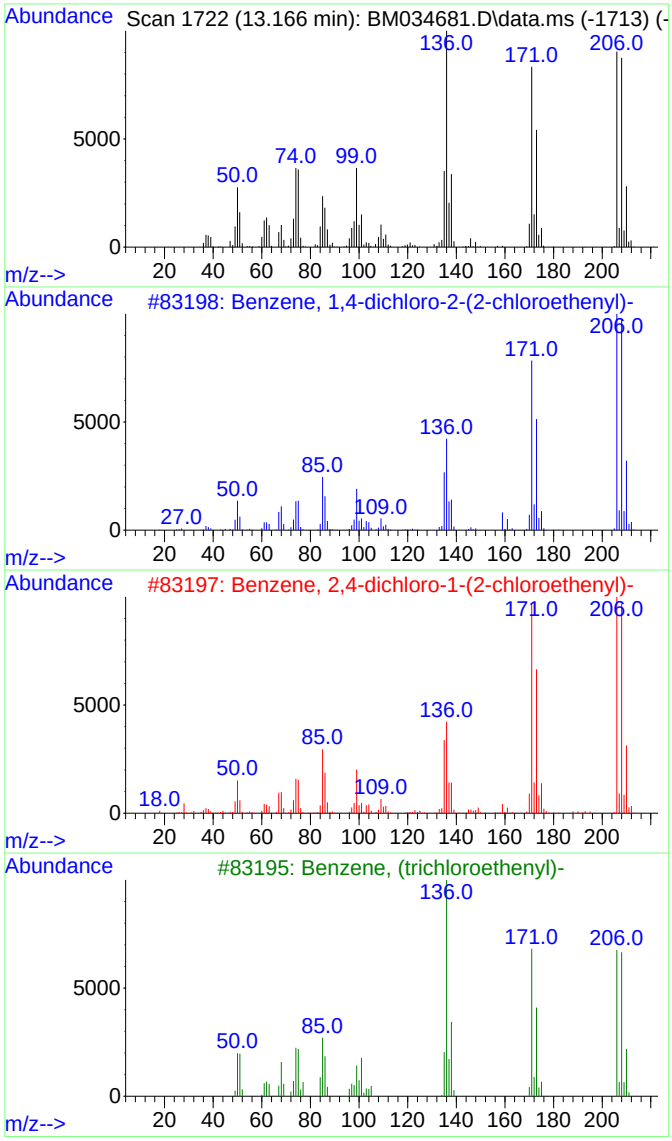
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

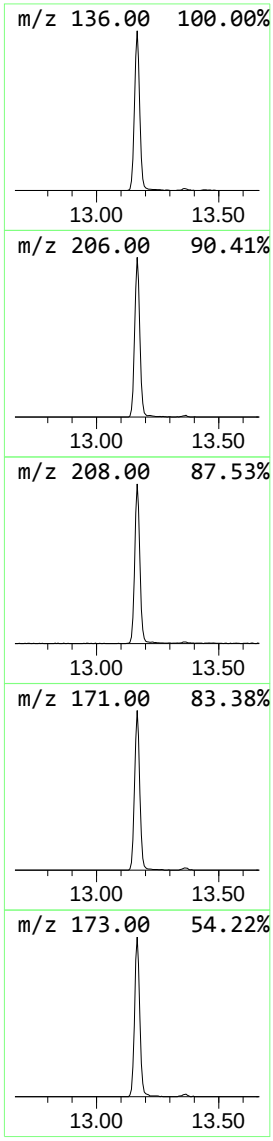
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1,4-dichloro-2-(2-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.166	18.55 ng/ul	1345890	Acenaphthene-d10	14.513

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-dichloro-2-(2-chlor...	206	C8H5Cl3	054935-00-1	96
2	Benzene, 2,4-dichloro-1-(2-chlor...	206	C8H5Cl3	045892-47-5	95
3	Benzene, (trichloroethenyl)-	206	C8H5Cl3	000700-60-7	90
4	Benzonitrile, 2,6-dichloro-	171	C7H3Cl2N	001194-65-6	55
5	1,1-Dichloro-2-(P-chlorophenyl)e...	206	C8H5Cl3	005263-17-2	53





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

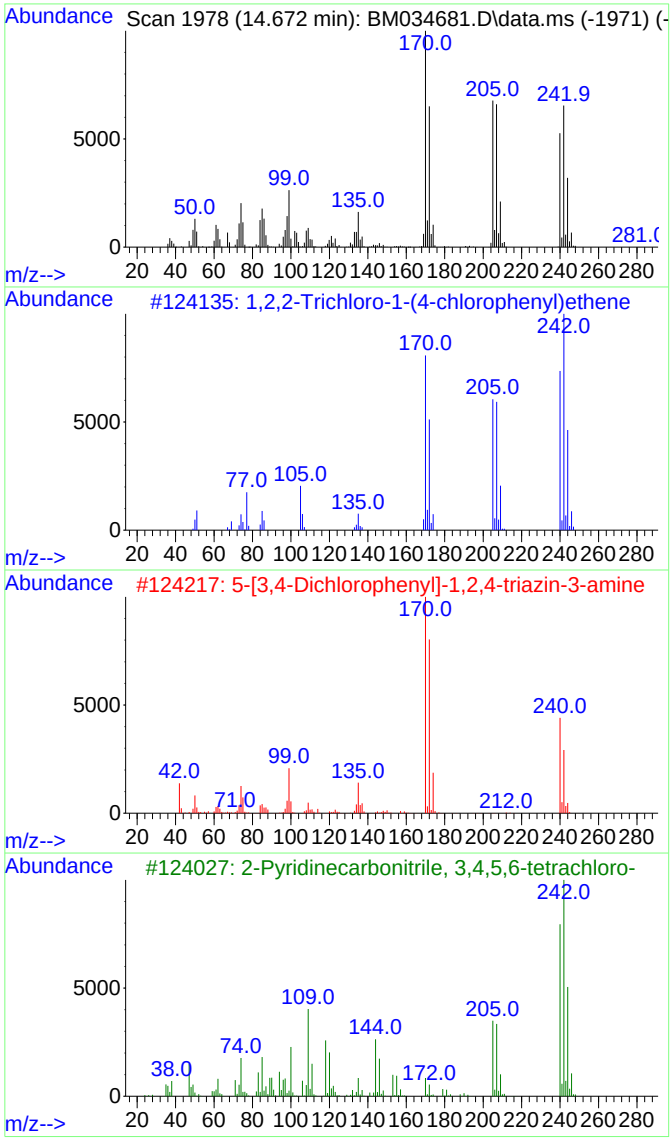
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

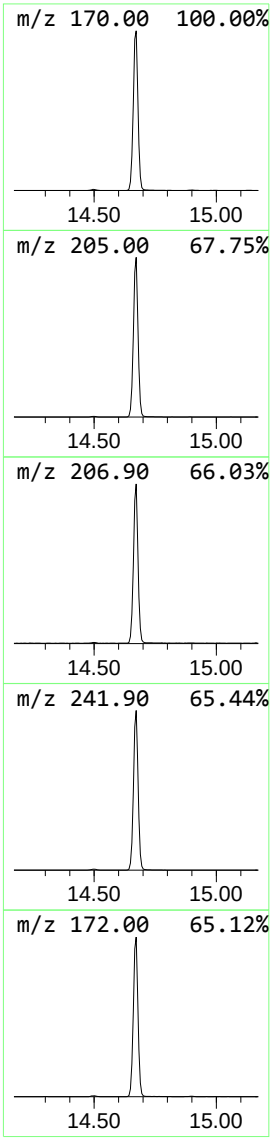
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1,2,2-Trichloro-1-(4-chloro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.672	69.74 ng/ul	5060030	Acenaphthene-d10	14.513

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,2-Trichloro-1-(4-chloropheny...	240	C8H4Cl4	004714-35-6	98
2	5-[3,4-Dichlorophenyl]-1,2,4-tri...	240	C9H6Cl2N4	1000213-66-2	49
3	2-Pyridinecarbonitrile, 3,4,5,6-...	240	C6Cl4N2	017824-83-8	35
4	2,4,6-Trichlorobenzonitrile	205	C7H2Cl3N	006575-05-9	35
5	5-Chloro-2-methoxy-2,4,6-cyclohe...	170	C8H7ClO2	013187-38-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

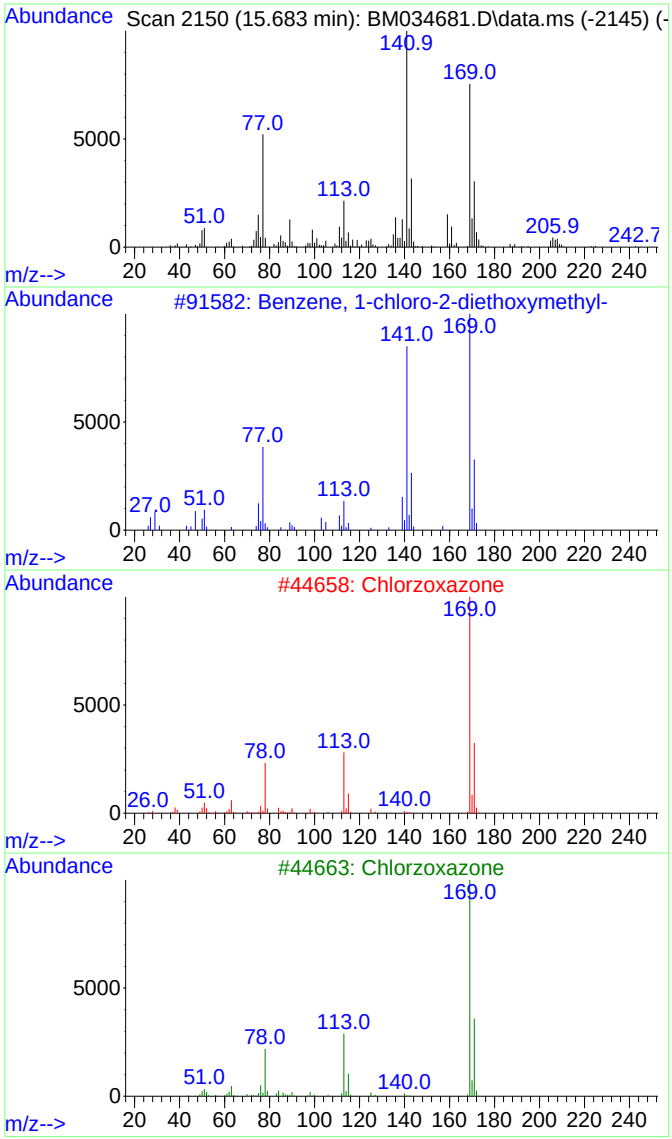
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

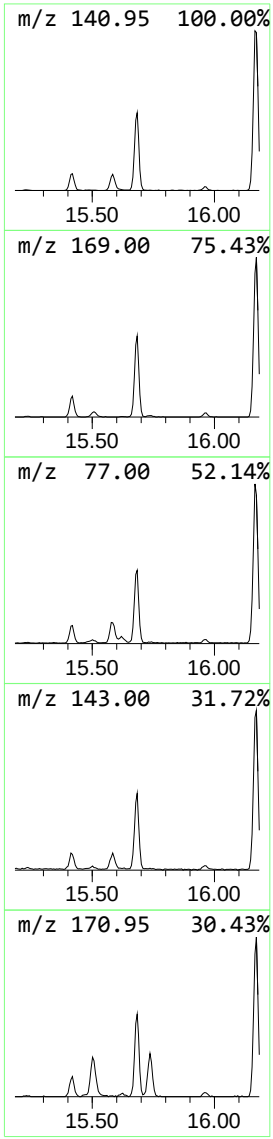
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 unknown-01 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.683	7.10 ng/ul	515388	Acenaphthene-d10	14.513

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-diethoxymethyl-	214	C11H15ClO2	035364-86-4	49
2	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	46
3	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	38
4	2-Chlorophenyl isothiocyanate	169	C7H4ClNS	002740-81-0	38
5	2-chloro-4,5-methylenedioxypheny...	212	C10H9ClO3	1000378-96-4	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

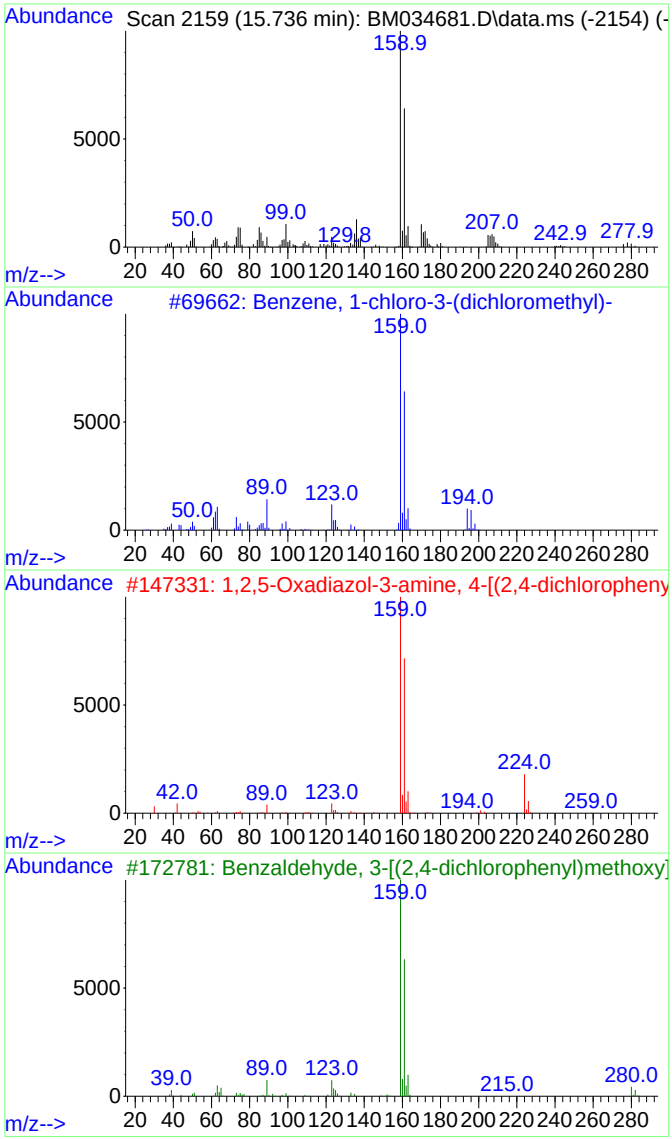
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

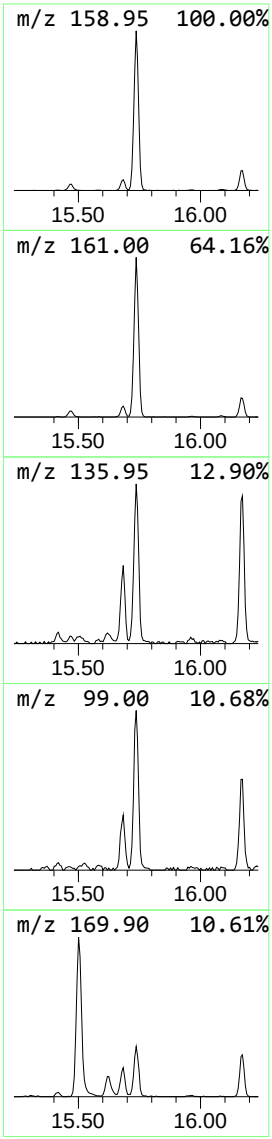
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-chloro-3-(dichlo... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.736	11.29 ng/ul	819040	Acenaphthene-d10	14.513

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-(dichloromet...	194	C7H5Cl3	015145-69-4	53
2	1,2,5-Oxadiazol-3-amine, 4-[(2,4...	259	C9H7Cl2N3O2	1000317-27-7	53
3	Benzaldehyde, 3-[(2,4-dichloroph...	280	C14H10Cl2O2	1000338-23-5	53
4	3,6-Bis[[[3,4-dichlorophenyl]met...	462	C16H10Cl4N4S2	075328-83-5	53
5	Benzene, 1,3-dichloro-2-ethyl-	174	C8H8Cl2	033407-02-2	53





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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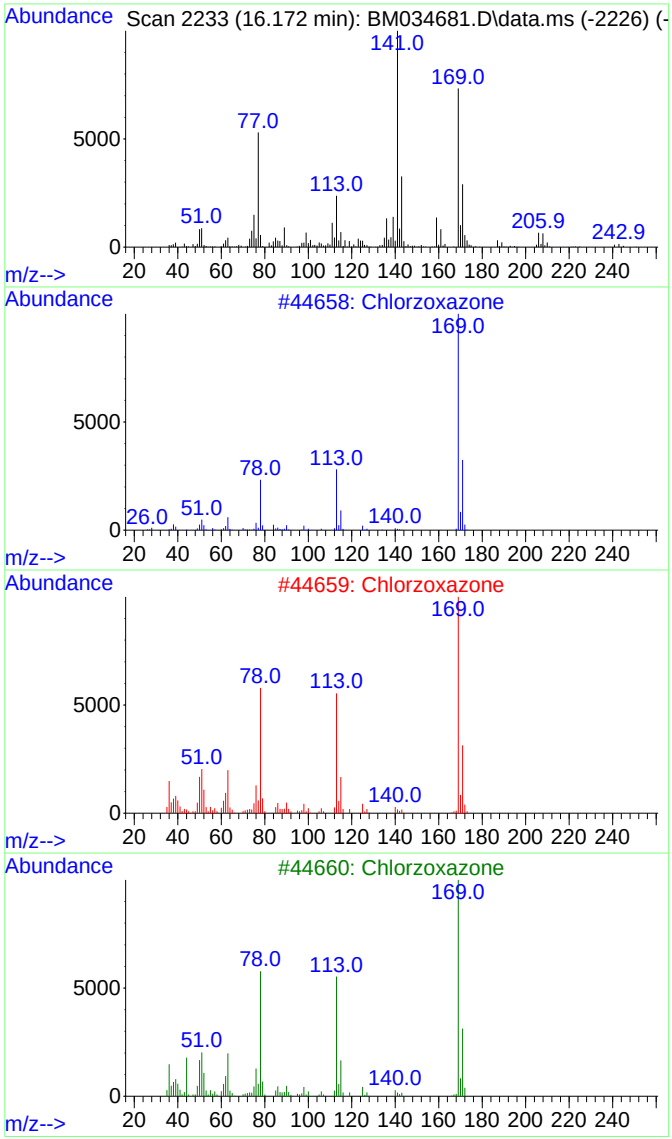
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

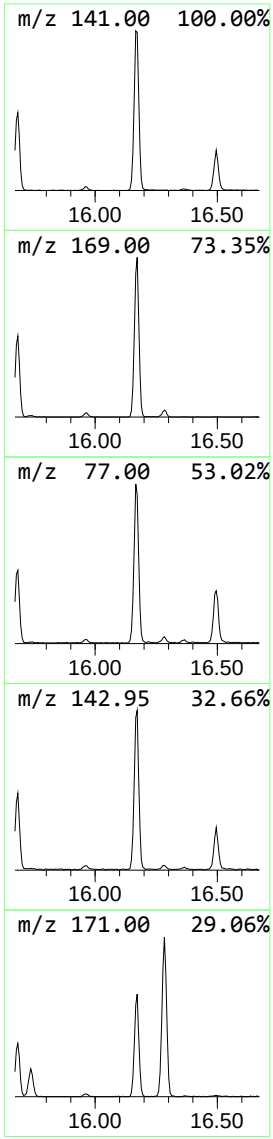
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown-02 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.172	13.83 ng/ul	1063010	Phenanthrene-d10	17.254

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	46
2	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	41
3	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	41
4	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	38
5	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	38





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ALS Vial : 23 Sample Multiplier: 1

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BNA_M
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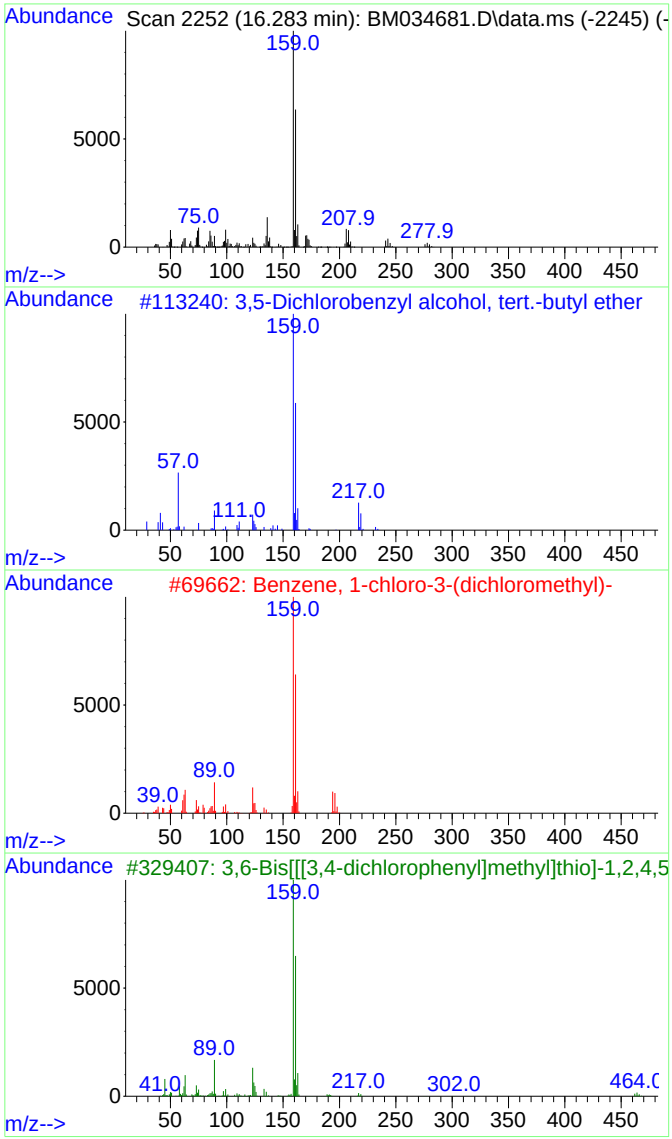
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

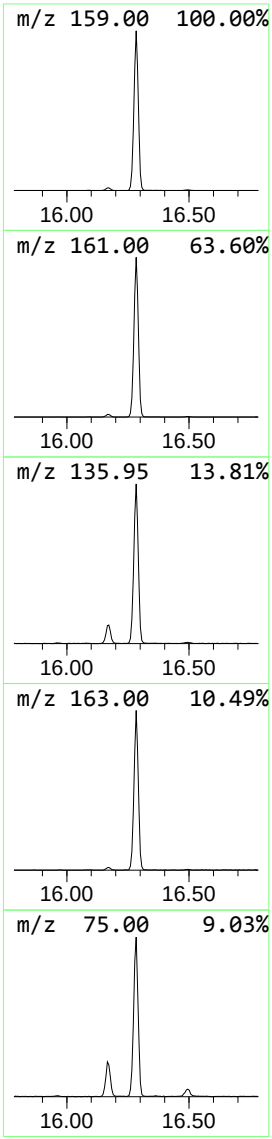
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3,5-Dichlorobenzyl alcohol,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.283	68.52 ng/ul	5267420	Phenanthrene-d10	17.254

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dichlorobenzyl alcohol, tert...	232	C11H14Cl2O	1000378-14-1	59
2	Benzene, 1-chloro-3-(dichloromet...	194	C7H5Cl3	015145-69-4	59
3	3,6-Bis[[[3,4-dichlorophenyl]met...	462	C16H10Cl4N4S2	075328-83-5	59
4	Benzene, (trichloromethyl)-	194	C7H5Cl3	000098-07-7	59
5	1,2,5-Oxadiazol-3-amine, 4-[(2,4...	259	C9H7Cl2N3O2	1000317-27-7	59





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

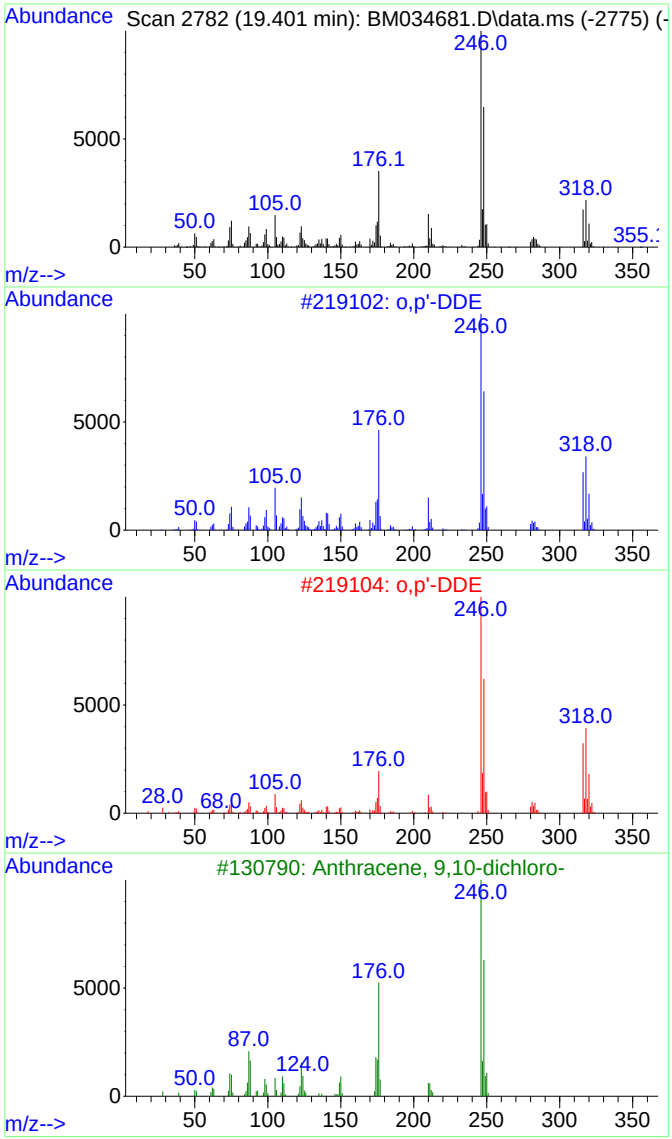
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

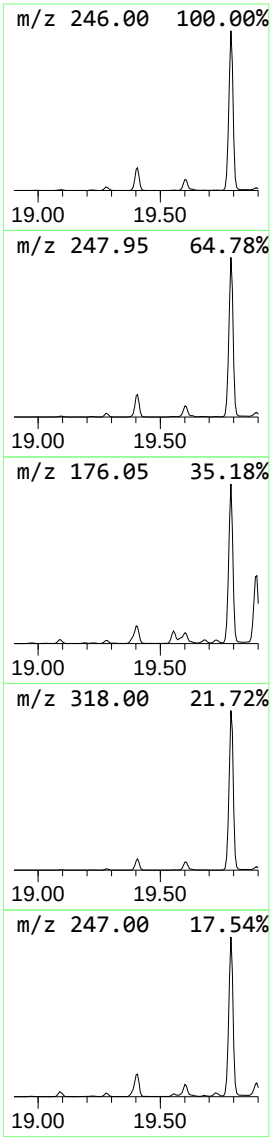
****

Peak Number 7 o,p'-DDE Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.401	24.70 ng/ul	1664210	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
2	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
3	Anthracene, 9,10-dichloro-	246	C14H8Cl2	000605-48-1	98
4	o,p'-DDE	316	C14H8Cl4	003424-82-6	96
5	p,p'-DDE	316	C14H8Cl4	000072-55-9	96





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

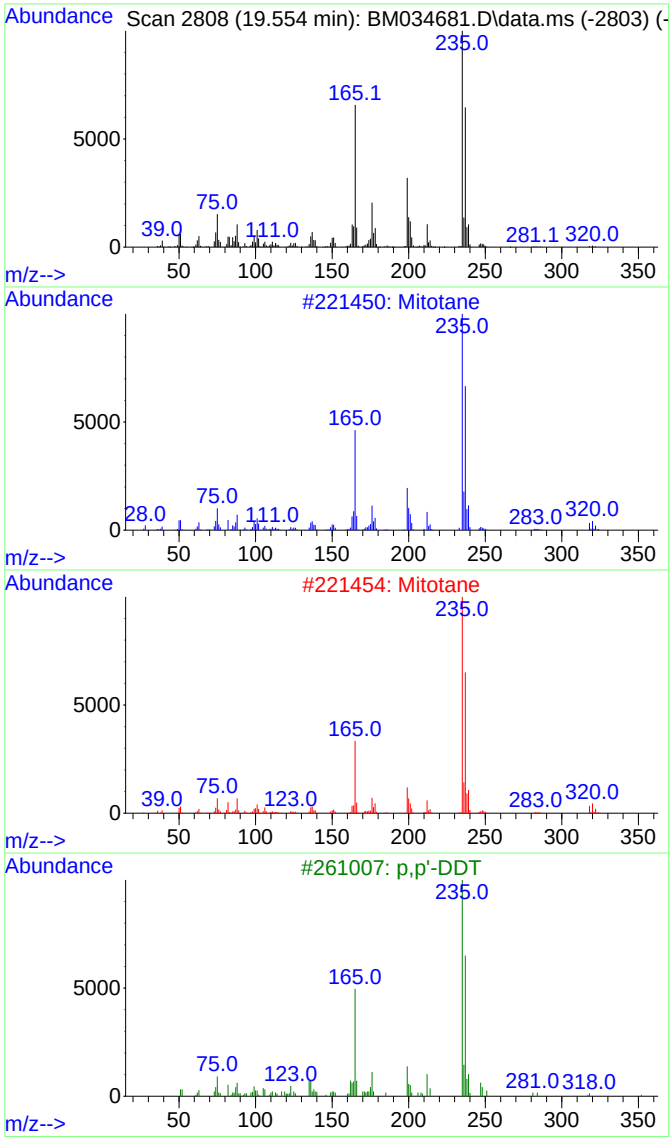
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

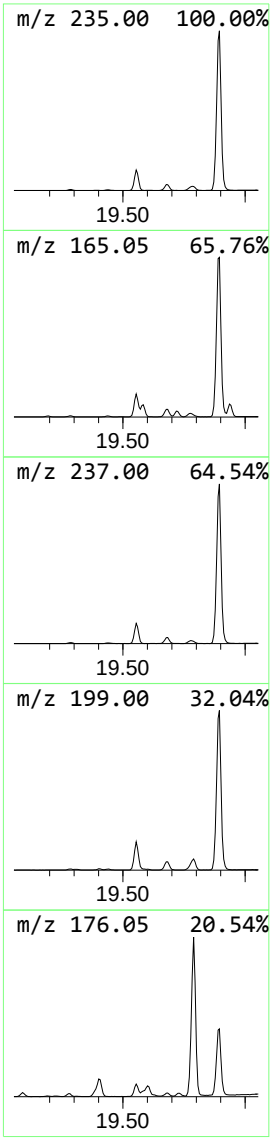
****

Peak Number 8 Mitotane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.554	25.00 ng/ul	1684590	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mitotane	318	C14H10Cl4	000053-19-0	91
2	Mitotane	318	C14H10Cl4	000053-19-0	87
3	p,p'-DDT	352	C14H9Cl5	000050-29-3	83
4	m,p'-DDD	318	C14H10Cl4	004329-12-8	83
5	m,p'-DDD	318	C14H10Cl4	004329-12-8	80





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

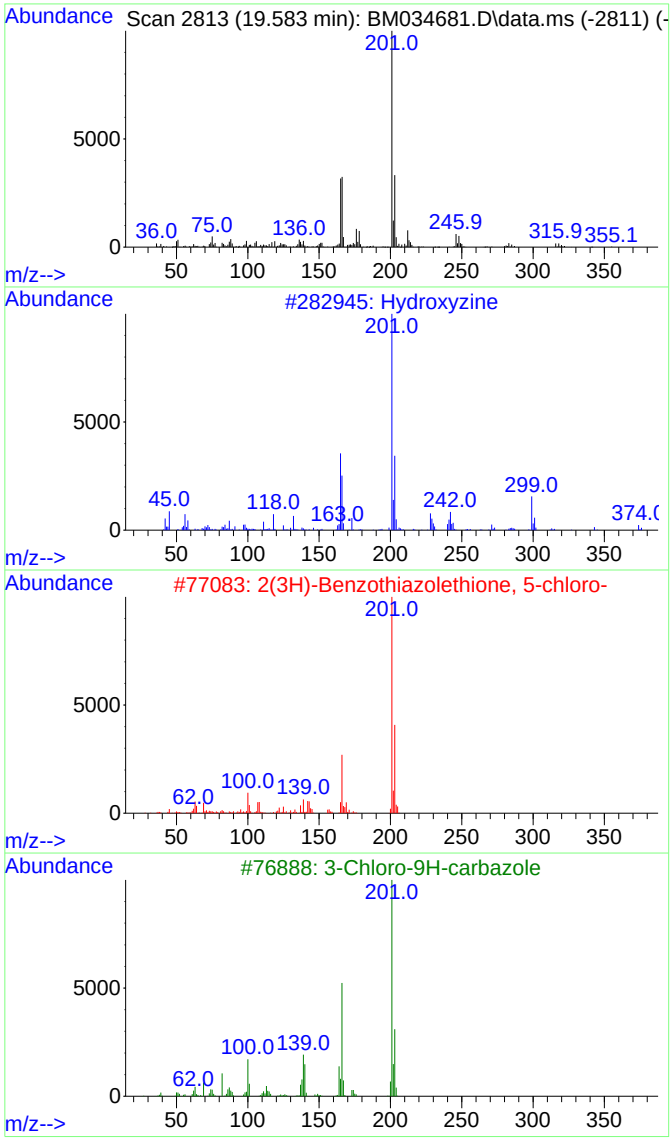
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

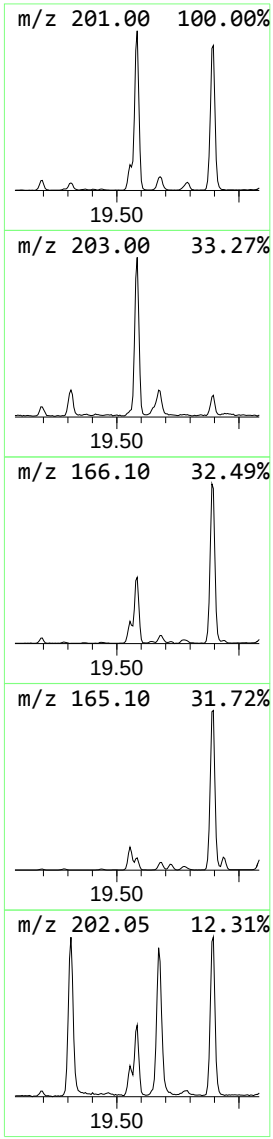
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Hydroxyzine Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.583	13.56 ng/ul	913708	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydroxyzine	374	C21H27ClN2O2	000068-88-2	72
2	2(3H)-Benzothiazolethione, 5-chl...	201	C7H4ClNS2	005331-91-9	64
3	3-Chloro-9H-carbazole	201	C12H8ClN	002732-25-4	64
4	Hydroxyzine	374	C21H27ClN2O2	000068-88-2	56
5	Acetic acid, (p-chlorophenyl)phe...	260	C15H13ClO2	005359-46-6	56





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Operator : CG/JU
Sample : N2316-13
Misc :
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Instrument :
BNA_M
ClientSampleId :
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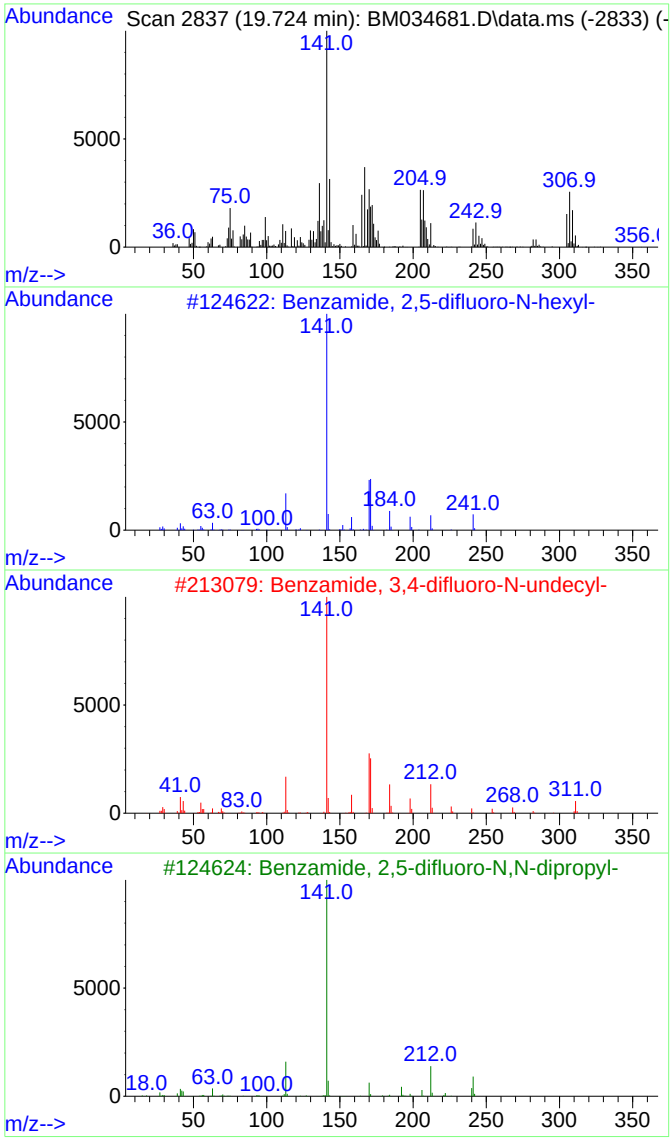
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

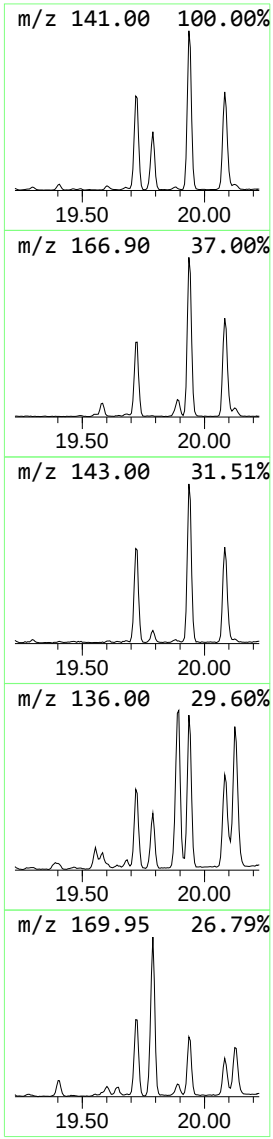
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown-03 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.724	22.27 ng/ul	1500860	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, 2,5-difluoro-N-hexyl-	241	C13H17F2NO	1010407-58-8	43
2	Benzamide, 3,4-difluoro-N-undecyl-	311	C18H27F2NO	1000407-80-0	43
3	Benzamide, 2,5-difluoro-N,N-dipr...	241	C13H17F2NO	1000438-25-7	38
4	Methyl 2,6-difluorobenzoate	172	C8H6F2O2	013671-00-6	35
5	Benzamide, 2,4-difluoro-N,N-dipr...	241	C13H17F2NO	1000438-27-3	35





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

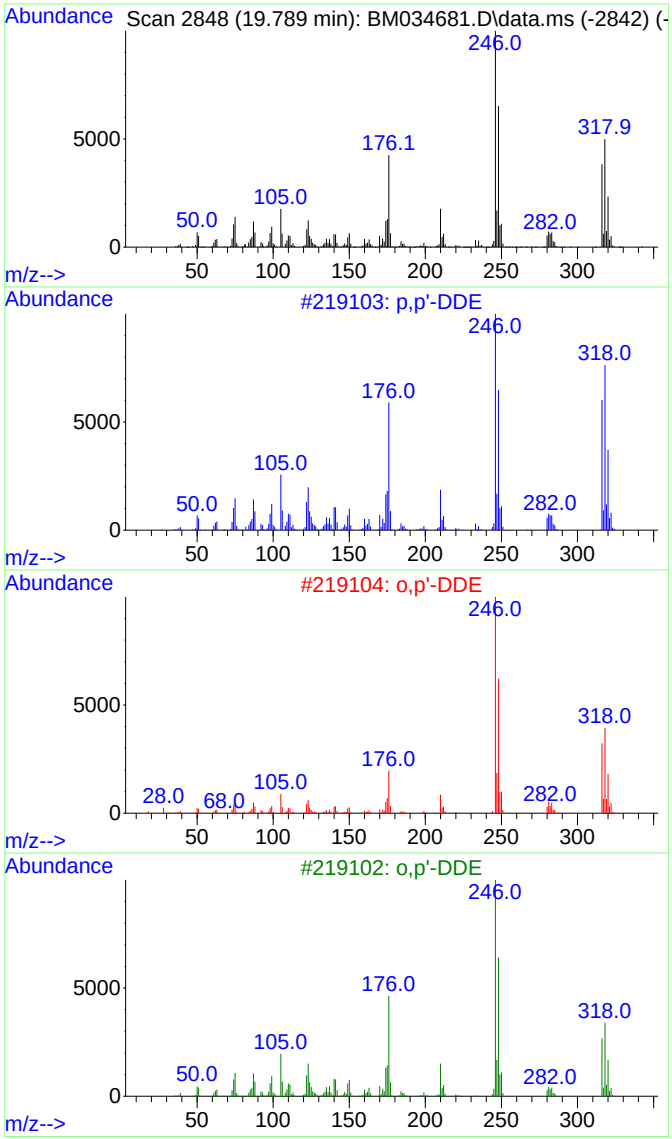
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

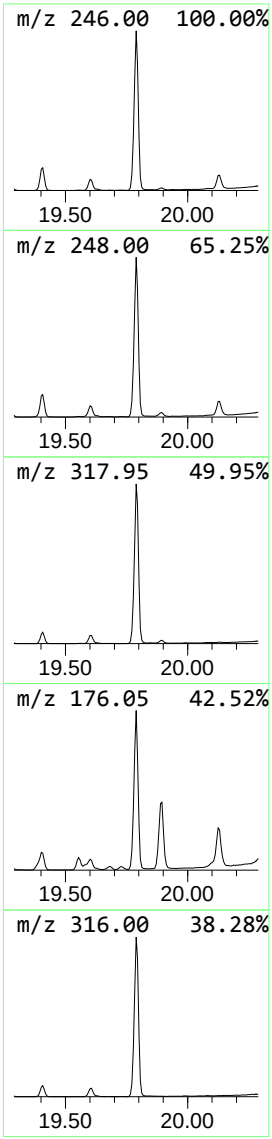
****

Peak Number 11 p,p'-DDE Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.789	180.36 ng/ul	12153100	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDE	316	C14H8Cl4	000072-55-9	99
2	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
3	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
4	p,p'-DDE	316	C14H8Cl4	000072-55-9	99
5	p,p'-DDE	316	C14H8Cl4	000072-55-9	99





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Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Title : SVOA CALIBRATION

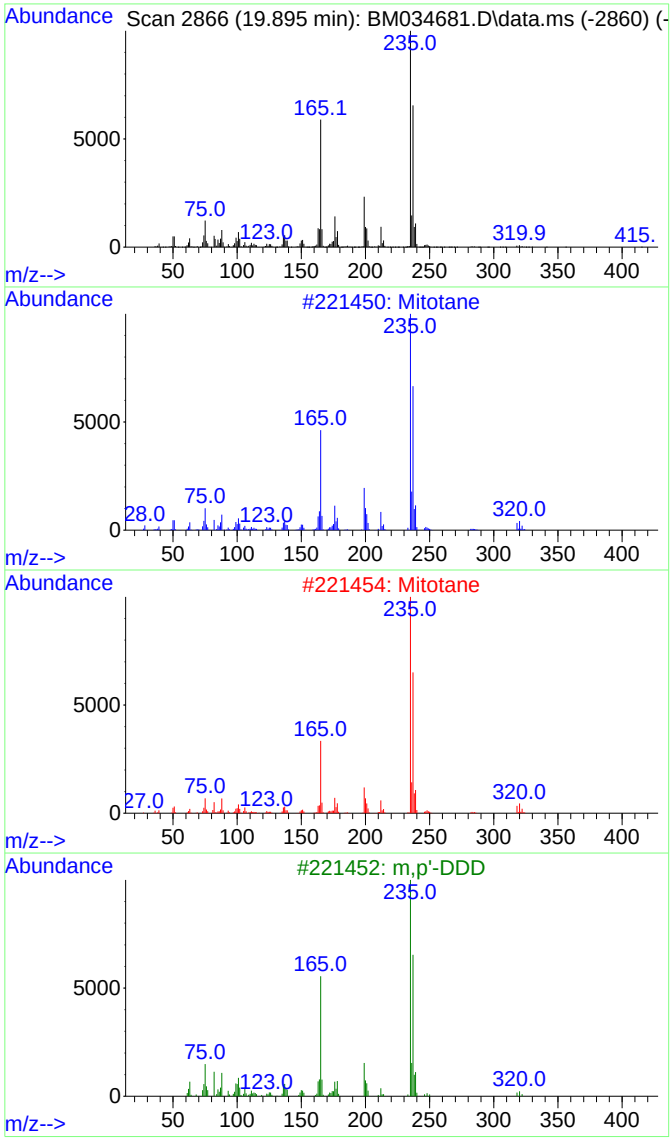
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

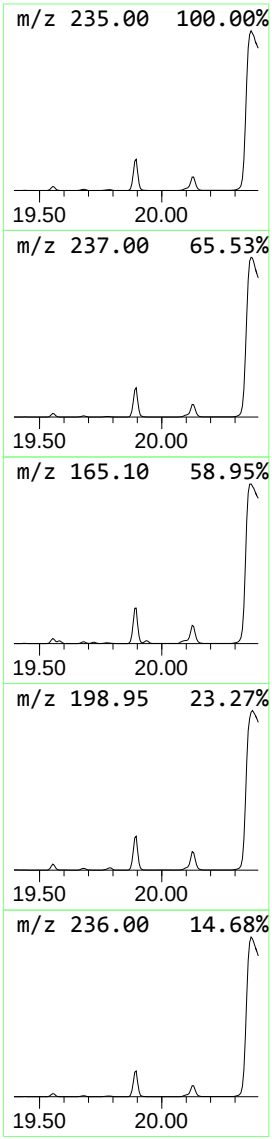
****

Peak Number 12 m,p'-DDD Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.895	176.68 ng/ul	11905500	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mitotane	318	C14H10Cl4	000053-19-0	96
2	Mitotane	318	C14H10Cl4	000053-19-0	91
3	m,p'-DDD	318	C14H10Cl4	004329-12-8	90
4	p,p'-DDT	352	C14H9Cl5	000050-29-3	90
5	m,p'-DDD	318	C14H10Cl4	004329-12-8	86





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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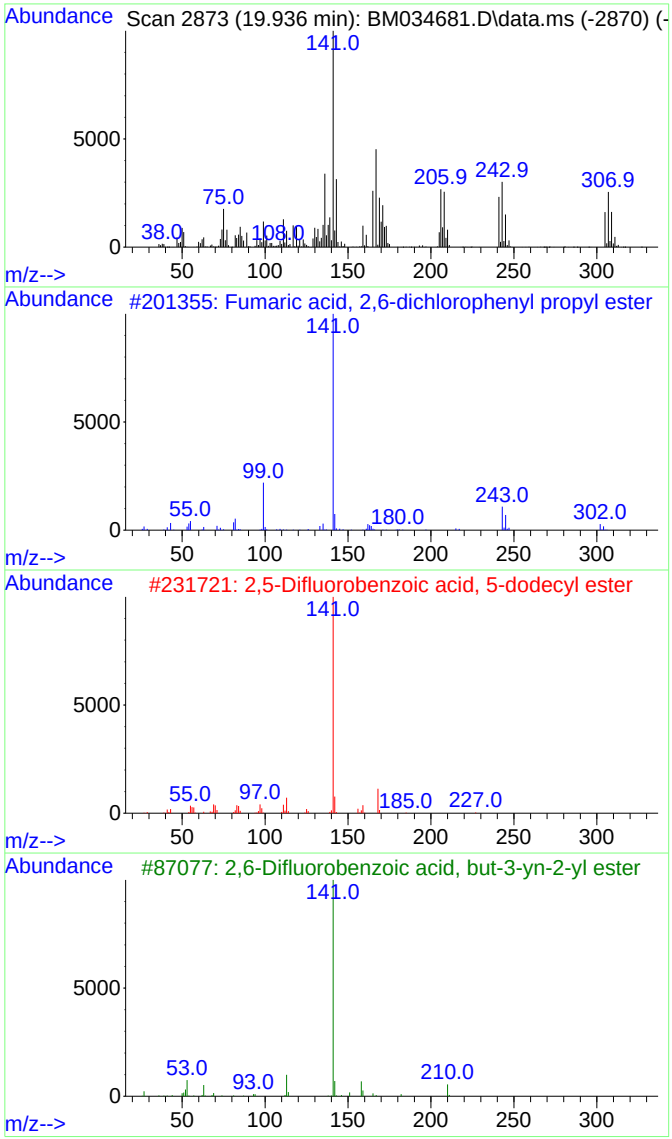
Quant Title : SVOA CALIBRATION

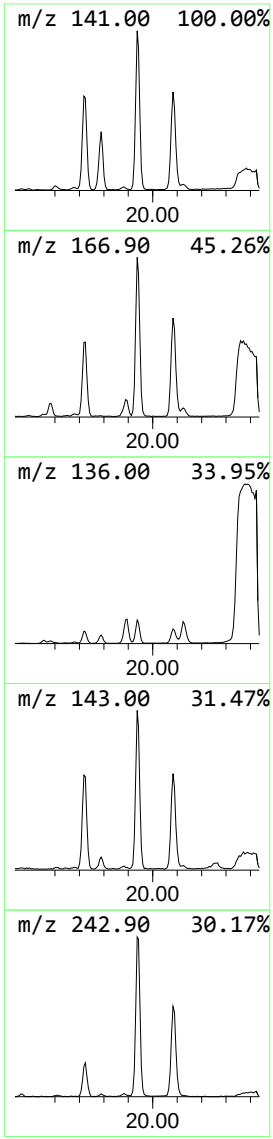
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 unknown-04 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.936	34.06 ng/ul	2294790	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fumaric acid, 2,6-dichlorophenyl...	302	C13H12Cl2O4	1000345-23-1	27
2	2,5-Difluorobenzoic acid, 5-dode...	326	C19H28F2O2	1000338-47-2	14
3	2,6-Difluorobenzoic acid, but-3-...	210	C11H8F2O2	1000292-58-1	14
4	2,4-Difluorobenzoic acid, 4-dode...	326	C19H28F2O2	1000338-56-4	14
5	2,4-Difluorobenzoic acid, 3-pent...	368	C22H34F2O2	1010338-57-7	14





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Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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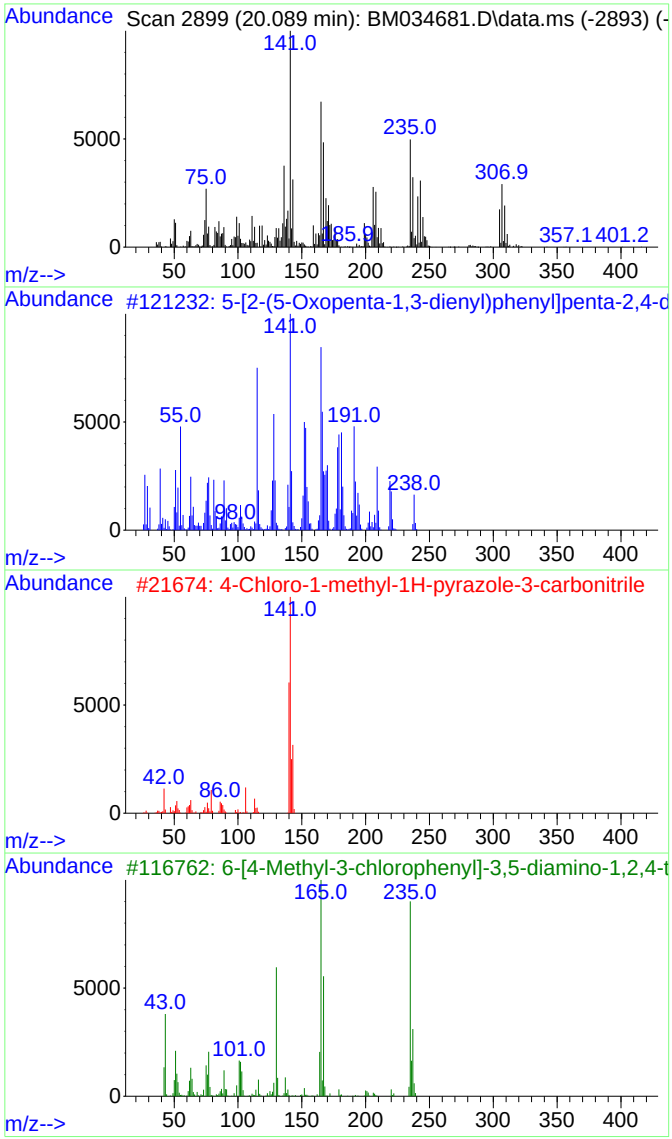
Quant Title : SVOA CALIBRATION

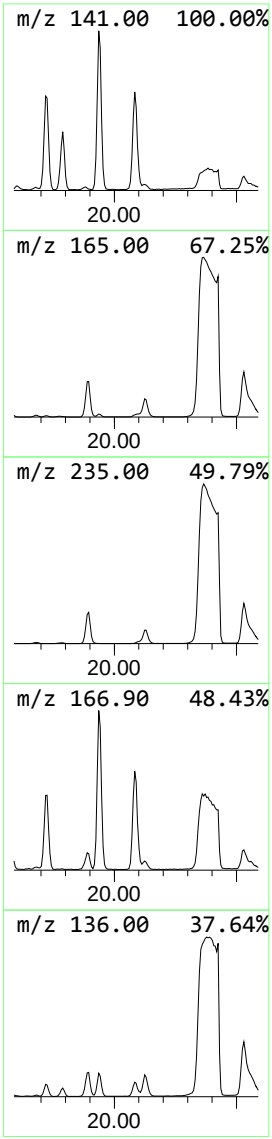
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 5-[2-(5-Oxopenta-1,3-dienyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.089	32.71 ng/ul	2203910	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-[2-(5-Oxopenta-1,3-dienyl)phen...	238	C16H14O2	1000192-91-9	74
2	4-Chloro-1-methyl-1H-pyrazole-3-...	141	C5H4ClN3	175204-86-1	25
3	6-[4-Methyl-3-chlorophenyl]-3,5-...	235	C10H10ClN5	058848-67-2	25
4	5-Amino-2-fluorophenol, methyl e...	141	C7H8FNO	064465-53-8	15
5	9H-Fluorene, 9-(1-methylethyl)-	208	C16H16	003299-99-8	15





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Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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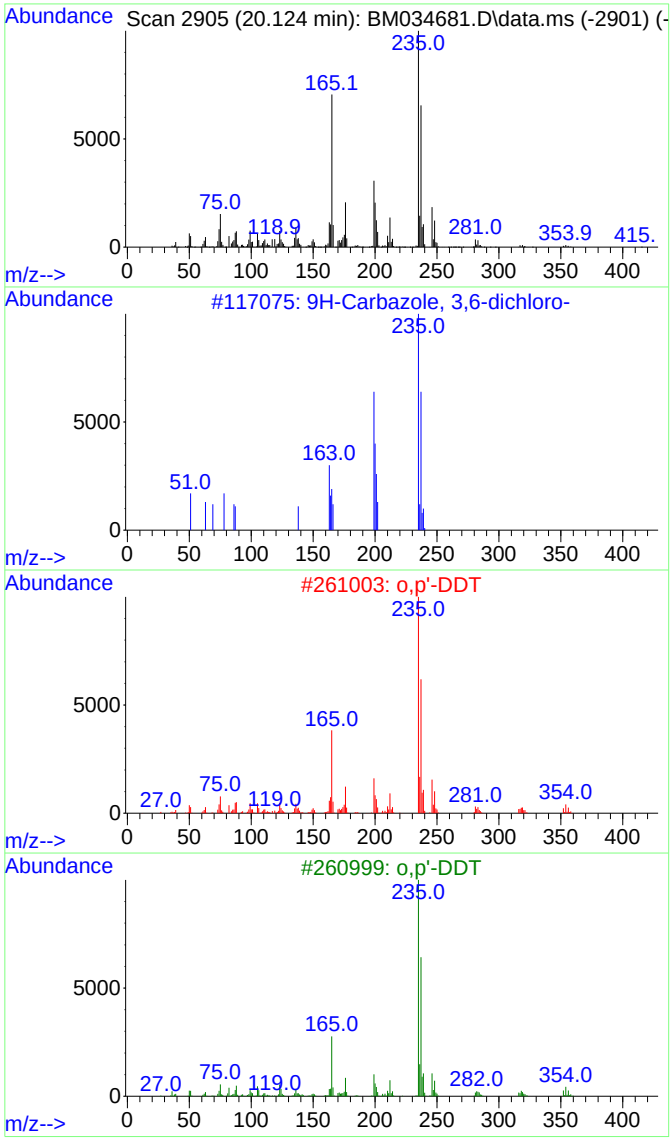
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

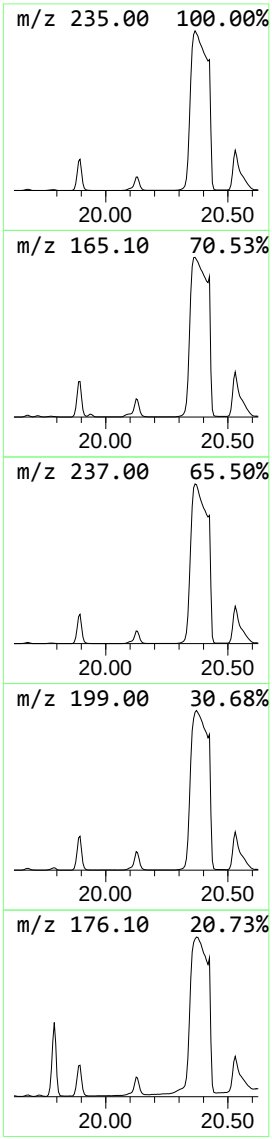
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 9H-Carbazole, 3,6-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.124	106.90 ng/ul	7203320	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 3,6-dichloro-	235	C12H7Cl2N	005599-71-3	94
2	o,p'-DDT	352	C14H9Cl5	000789-02-6	94
3	o,p'-DDT	352	C14H9Cl5	000789-02-6	93
4	o,p'-DDT	352	C14H9Cl5	000789-02-6	90
5	Mitotane	318	C14H10Cl4	000053-19-0	72





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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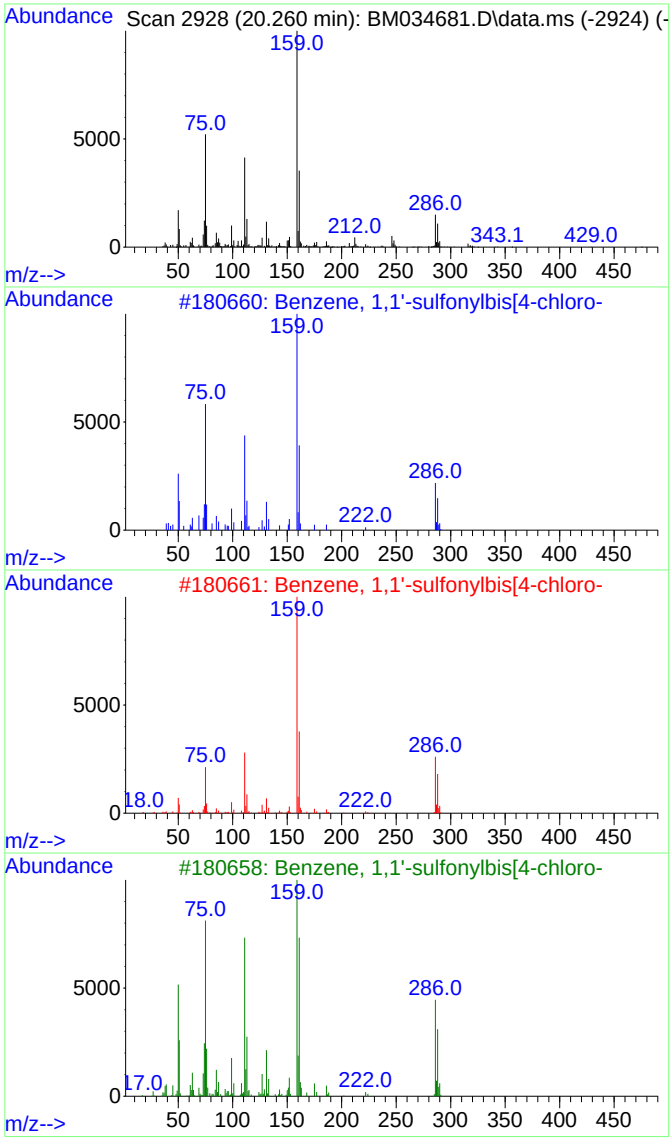
Quant Title : SVOA CALIBRATION

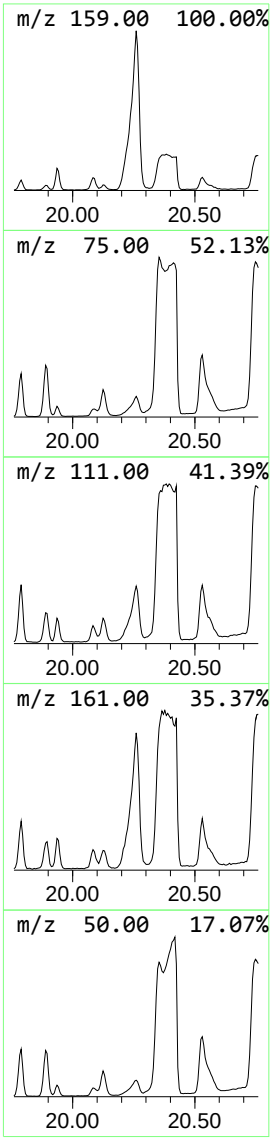
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1,1'-sulfonylbis[4... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.260	10.08 ng/ul	679339	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-sulfonylbis[4-chloro-	286	C12H8Cl2O2S	000080-07-9	99
2	Benzene, 1,1'-sulfonylbis[4-chloro-	286	C12H8Cl2O2S	000080-07-9	98
3	Benzene, 1,1'-sulfonylbis[4-chloro-	286	C12H8Cl2O2S	000080-07-9	97
4	Benzene, 1,1'-sulfonylbis[4-chloro-	286	C12H8Cl2O2S	000080-07-9	95
5	Benzene, 1,1'-sulfonylbis[4-chloro-	286	C12H8Cl2O2S	000080-07-9	91





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Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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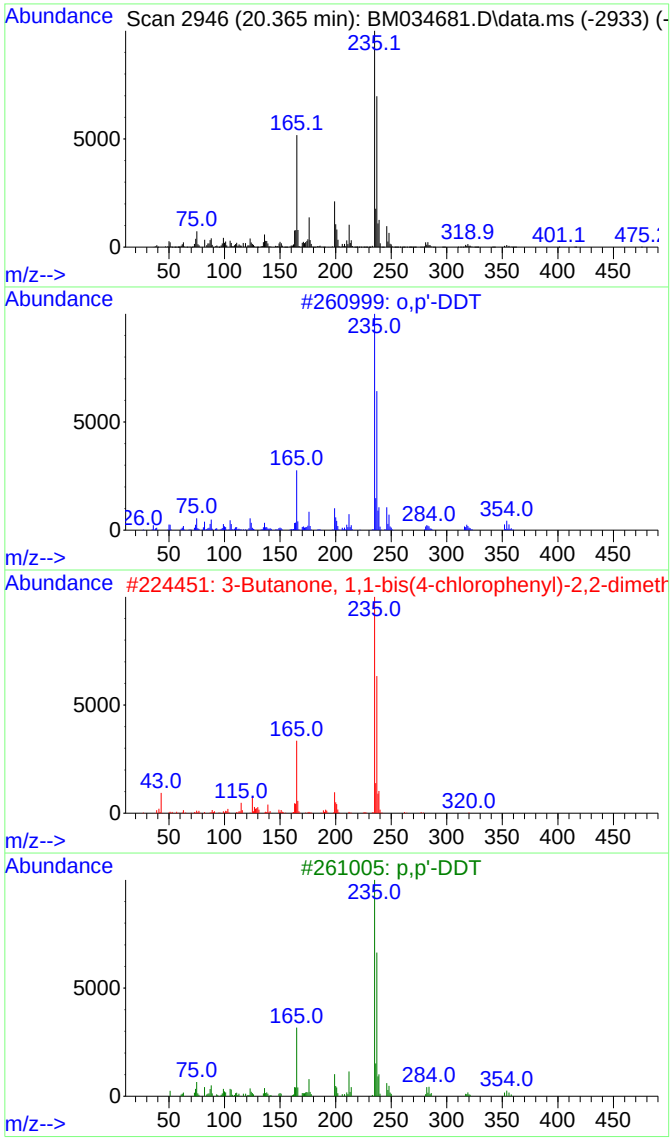
Quant Title : SVOA CALIBRATION

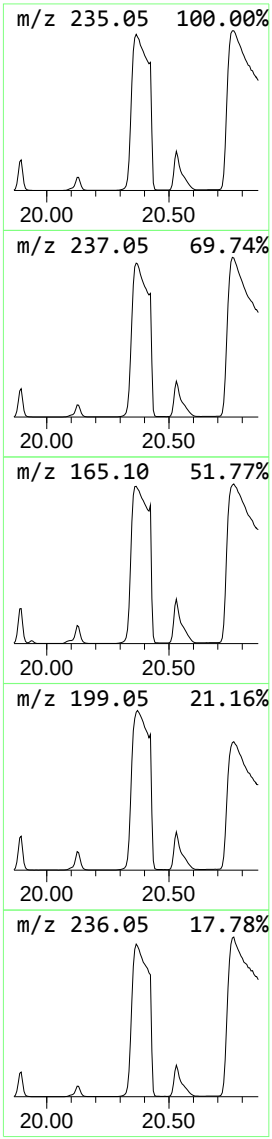
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 o,p'-DDT Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.366	3022.36 ng/ul	203659000	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDT	352	C14H9Cl5	000789-02-6	97
2	3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	93
3	p,p'-DDT	352	C14H9Cl5	000050-29-3	93
4	Mitotane	318	C14H10Cl4	000053-19-0	91
5	Mitotane	318	C14H10Cl4	000053-19-0	91





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

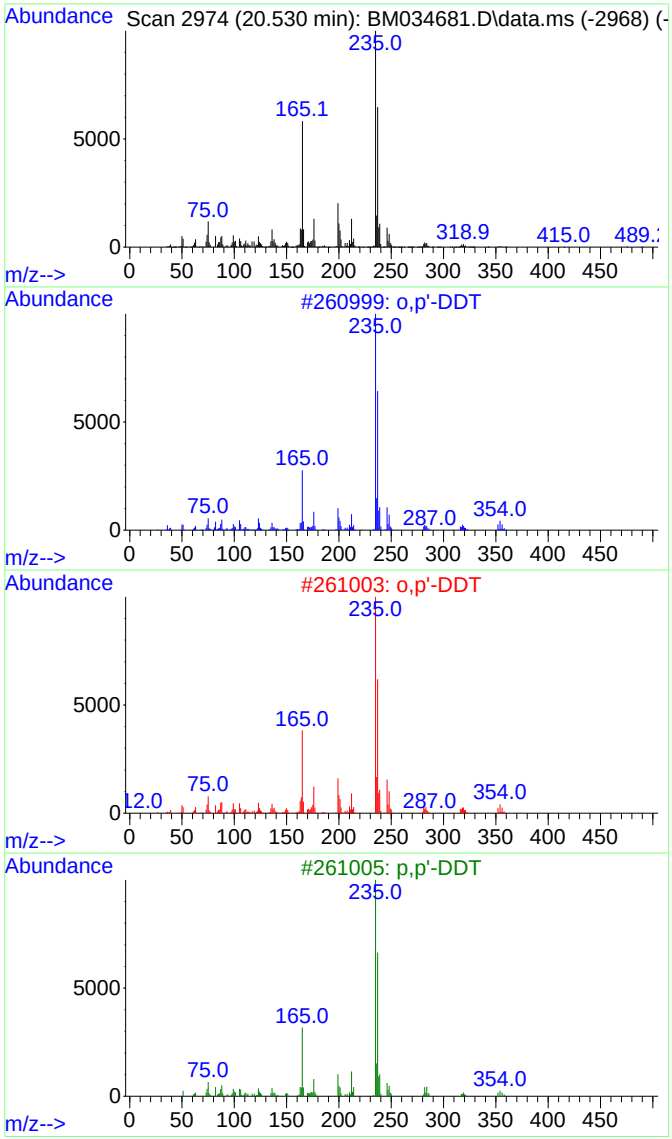
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

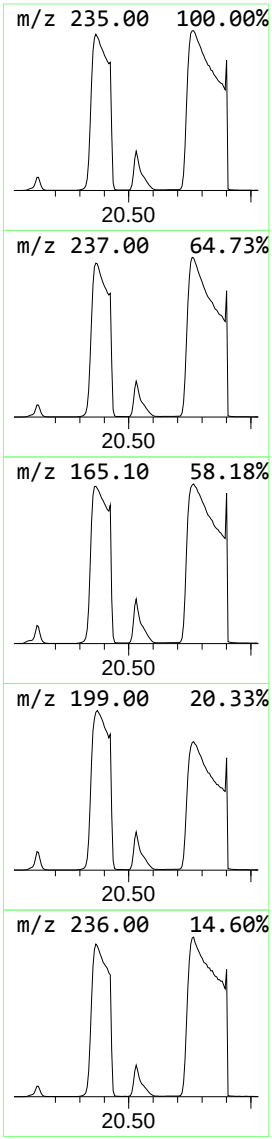
****

Peak Number 18 p,p'-DDT Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.530	406.43 ng/ul	27387000	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDT	352	C14H9Cl5	000789-02-6	98
2	o,p'-DDT	352	C14H9Cl5	000789-02-6	93
3	p,p'-DDT	352	C14H9Cl5	000050-29-3	91
4	o,p'-DDT	352	C14H9Cl5	000789-02-6	91
5	Mitotane	318	C14H10Cl4	000053-19-0	91





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ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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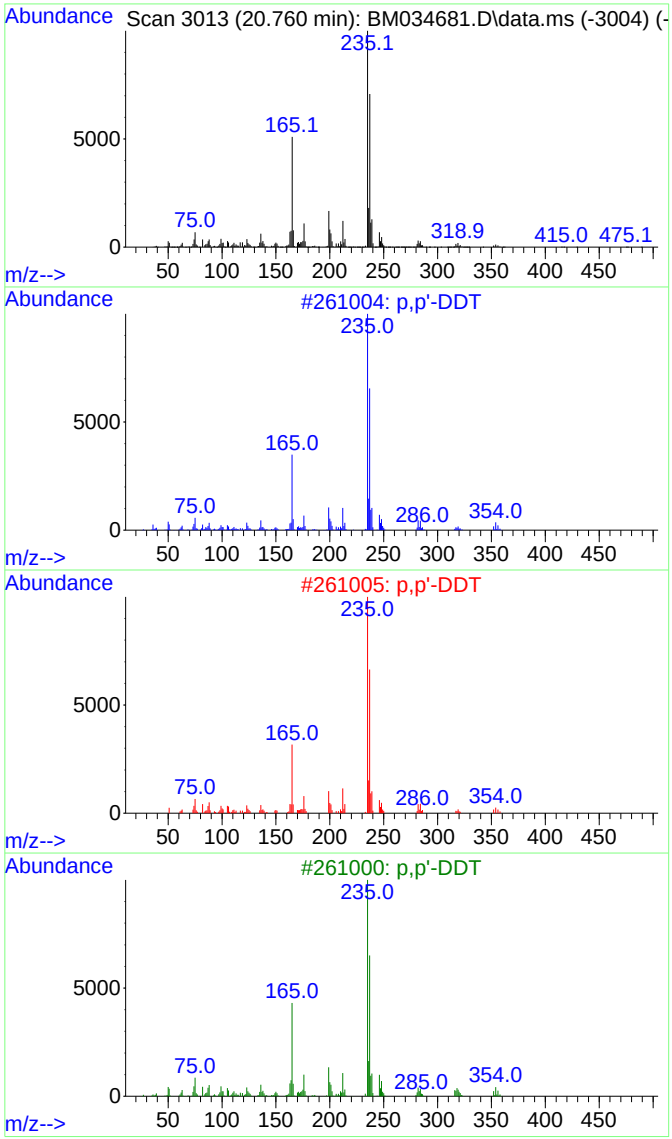
Quant Title : SVOA CALIBRATION

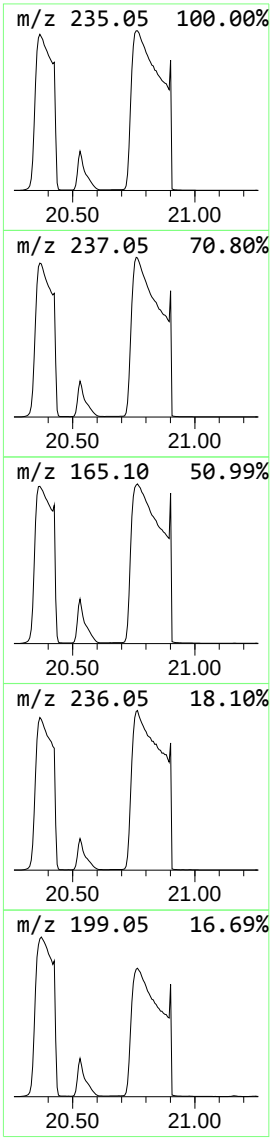
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 unknown-05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.760	5433.88 ng/ul	366156000	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDT	352	C14H9Cl5	000050-29-3	99
2	p,p'-DDT	352	C14H9Cl5	000050-29-3	98
3	p,p'-DDT	352	C14H9Cl5	000050-29-3	94
4	p,p'-DDT	352	C14H9Cl5	000050-29-3	94
5	o,p'-DDT	352	C14H9Cl5	000789-02-6	94





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

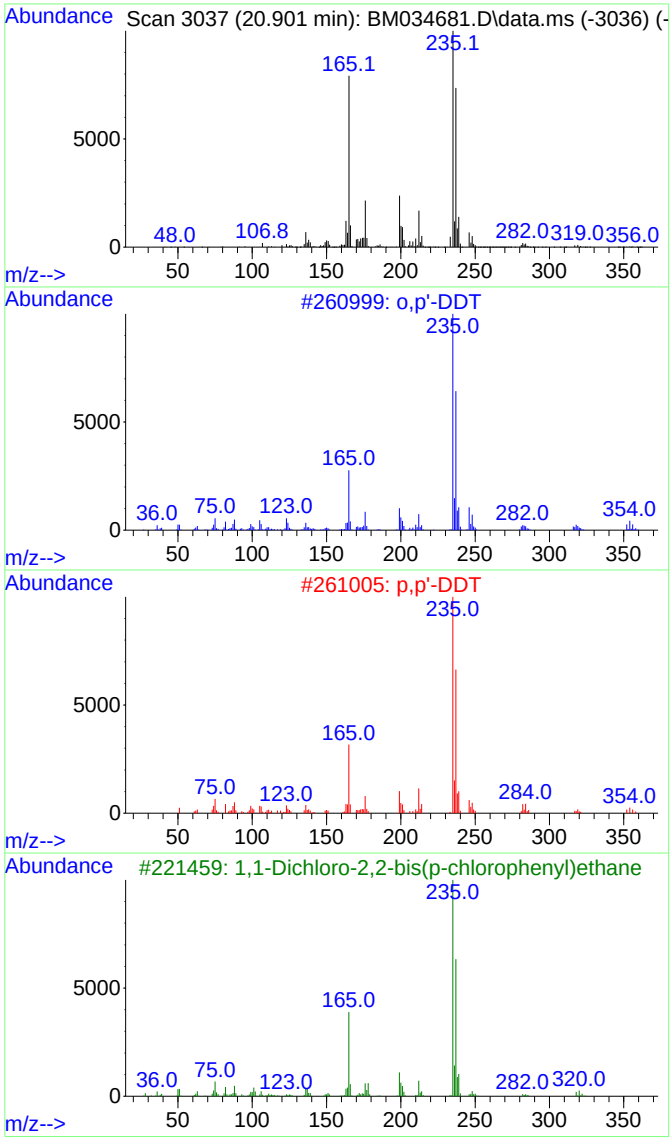
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

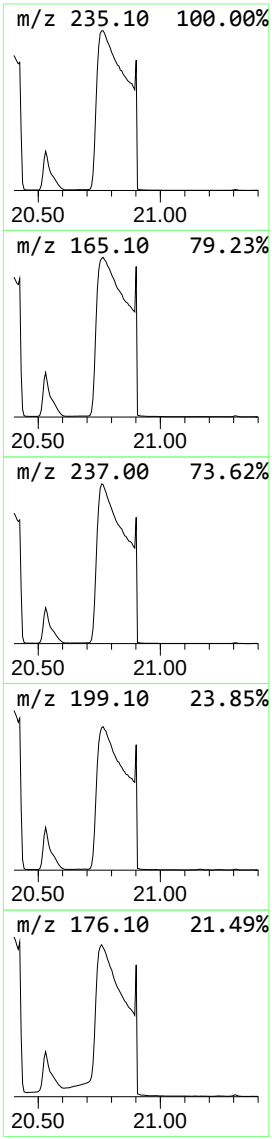
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 1,1-Dichloro-2,2-bis(p-chlo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.901	188.70 ng/ul	12715300	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDT	352	C14H9Cl5	000789-02-6	95
2	p,p'-DDT	352	C14H9Cl5	000050-29-3	89
3	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	86
4	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	86
5	1-Chloro-2,2-Bis(p-chlorophenyl)...	284	C14H11Cl3	002642-80-0	80





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

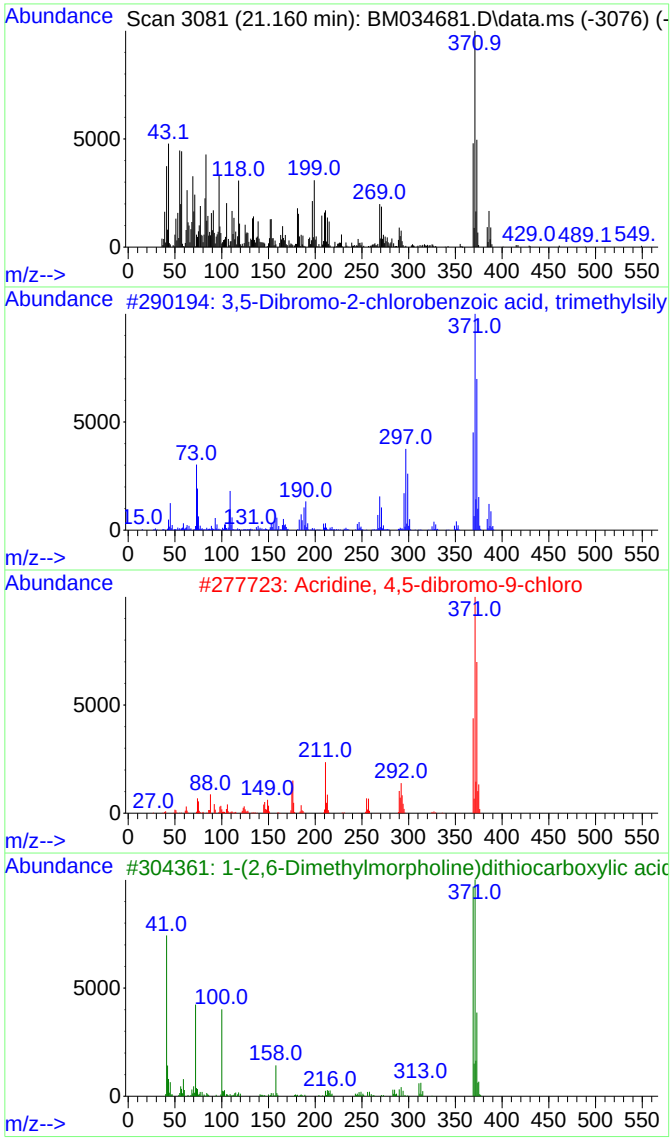
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

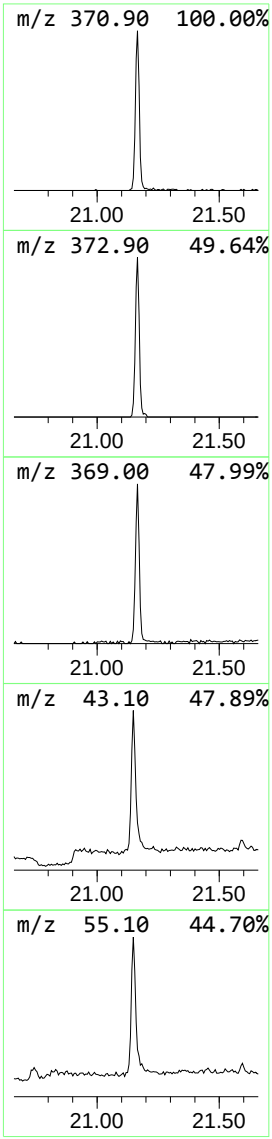
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 3,5-Dibromo-2-chlorobenzoic... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.160	12.29 ng/ul	828059	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dibromo-2-chlorobenzoic acid...	384	C10H11Br2ClO2Si	1000471-63-7	64
2	Acridine, 4,5-dibromo-9-chloro	369	C13H6Br2ClN	209460-13-9	38
3	1-(2,6-Dimethylmorpholine)dithio...	404	C12H12Cl4N2OS2	1000305-22-8	37
4	2-(4'-Trimethylsilyloxyphenyl)-2...	386	C22H34O2Si2	1000283-55-8	35
5	1-Amino-4-anilinoanthra-9,10-qui...	386	C23H22N2O2Si	1000474-79-9	30





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

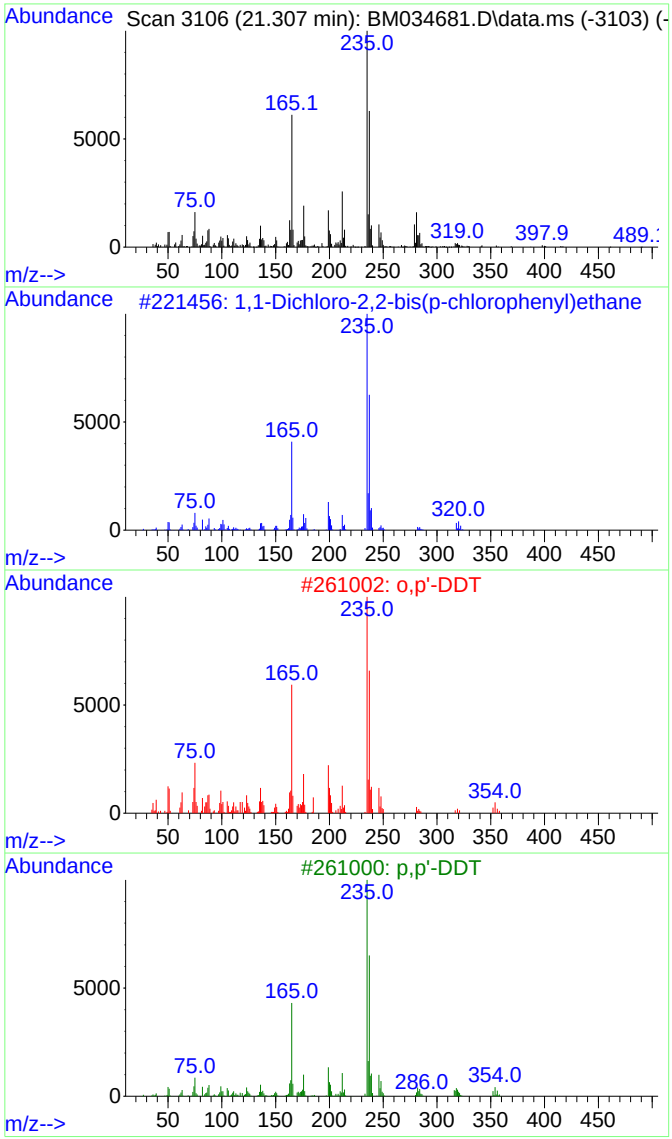
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

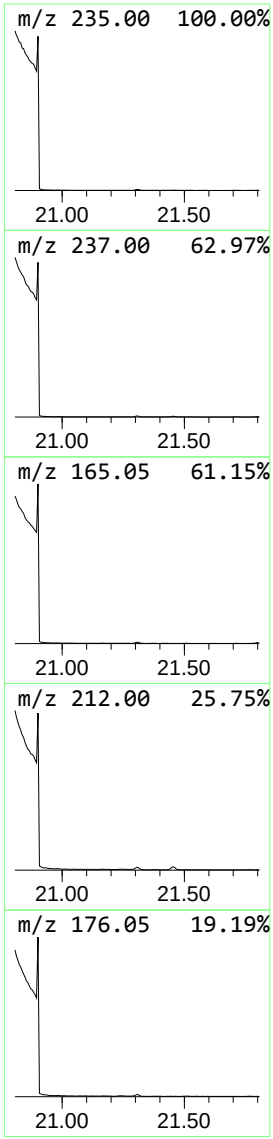
****

Peak Number 22 unknown-06 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.307	5.82 ng/ul	391892	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	86
2	o,p'-DDT	352	C14H9Cl5	000789-02-6	80
3	p,p'-DDT	352	C14H9Cl5	000050-29-3	74
4	p,p'-DDT	352	C14H9Cl5	000050-29-3	74
5	p,p'-DDT	352	C14H9Cl5	000050-29-3	70





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Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

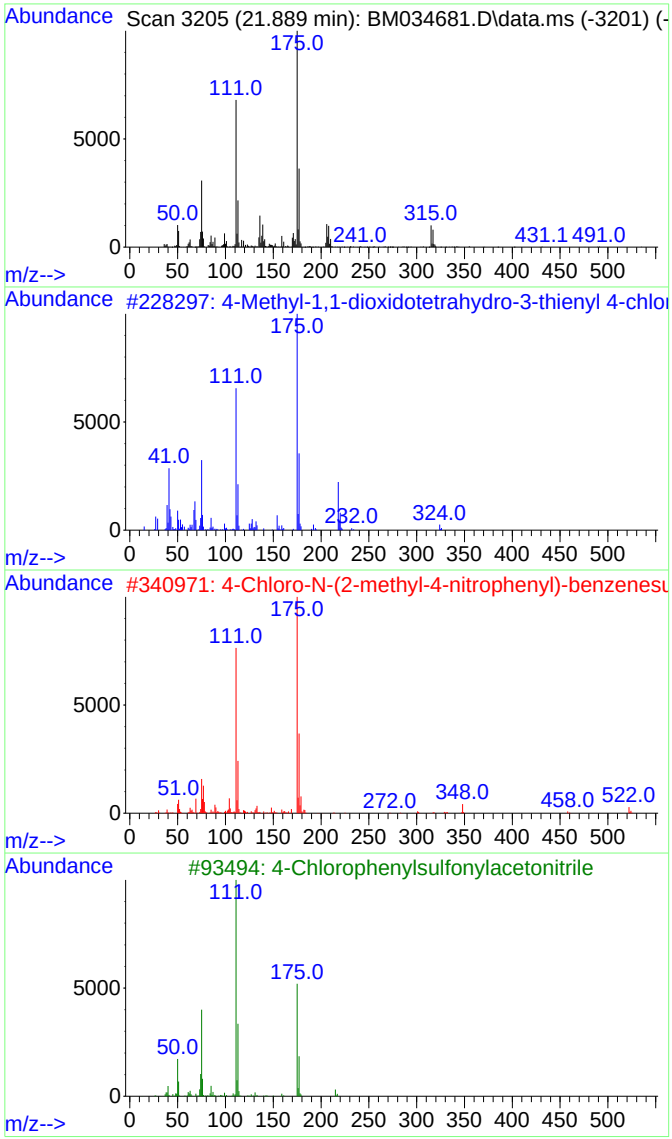
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

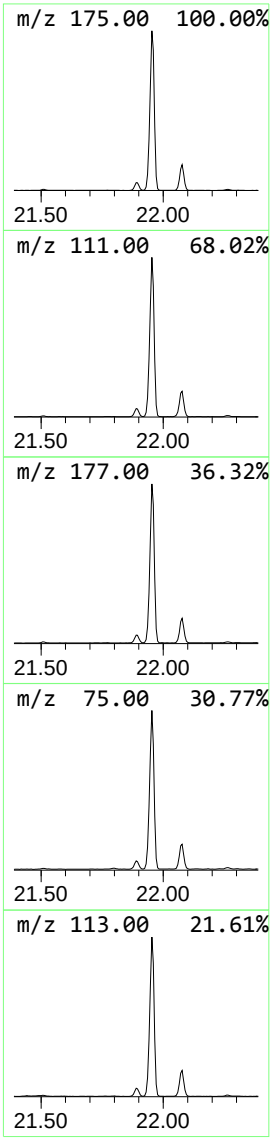
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 4-Methyl-1,1-dioxidotetrahy... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.889	10.79 ng/ul	726789	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-1,1-dioxidotetrahydro-3...	324	C11H13ClO5S2	1000332-40-7	59
2	4-Chloro-N-(2-methyl-4-nitrophen...	522	C17H10ClF7N2O5S	1000374-79-0	59
3	4-Chlorophenylsulfonylacetonitrile	215	C8H6ClNO2S	001851-09-8	50
4	Chloropropamide	276	C10H13ClN2O3S	000094-20-2	45
5	Benzenesulfonyl chloride, 4-chloro-	210	C6H4Cl2O2S	000098-60-2	45





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

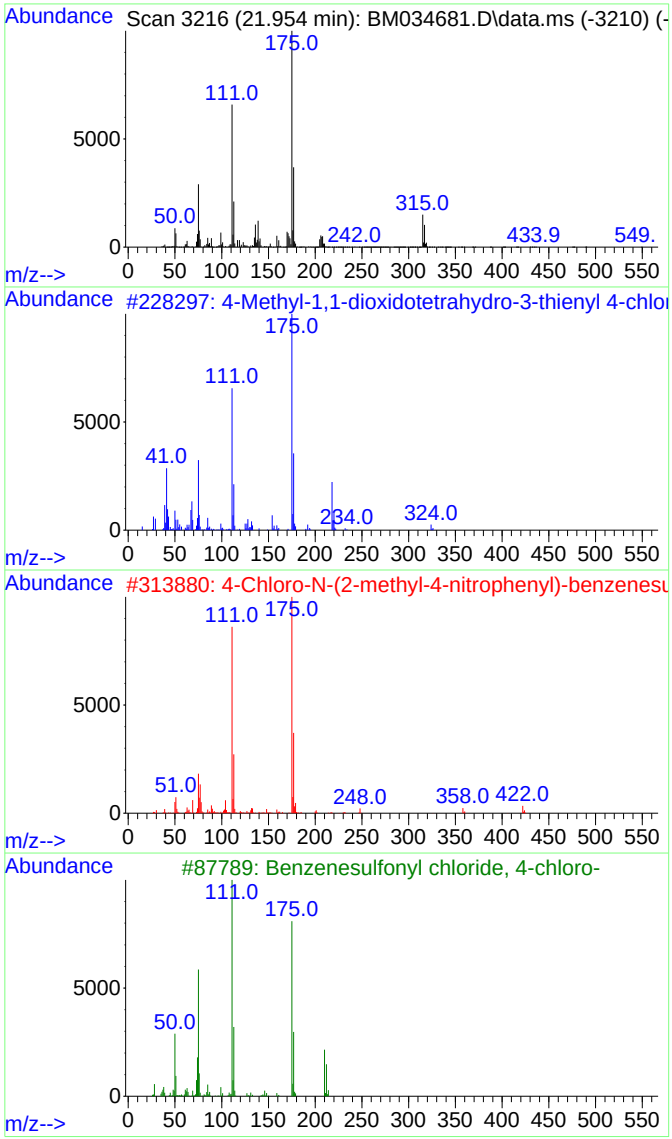
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

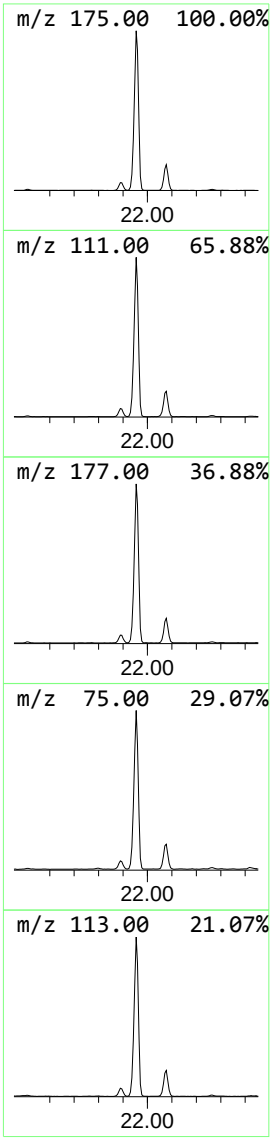
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 4-Chloro-N-(2-methyl-4-nitr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.954	197.02 ng/ul	13276000	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-1,1-dioxidotetrahydro-3...	324	C11H13ClO5S2	1000332-40-7	72
2	4-Chloro-N-(2-methyl-4-nitrophen...	422	C15H10ClF3N2O5S	1000374-78-8	64
3	Benzenesulfonyl chloride, 4-chloro-	210	C6H4Cl2O2S	000098-60-2	64
4	4-Chlorobenzenesulfonyl isocyanate	217	C7H4ClNO3S	005769-15-3	50
5	Chloropropamide	276	C10H13ClN2O3S	000094-20-2	50





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

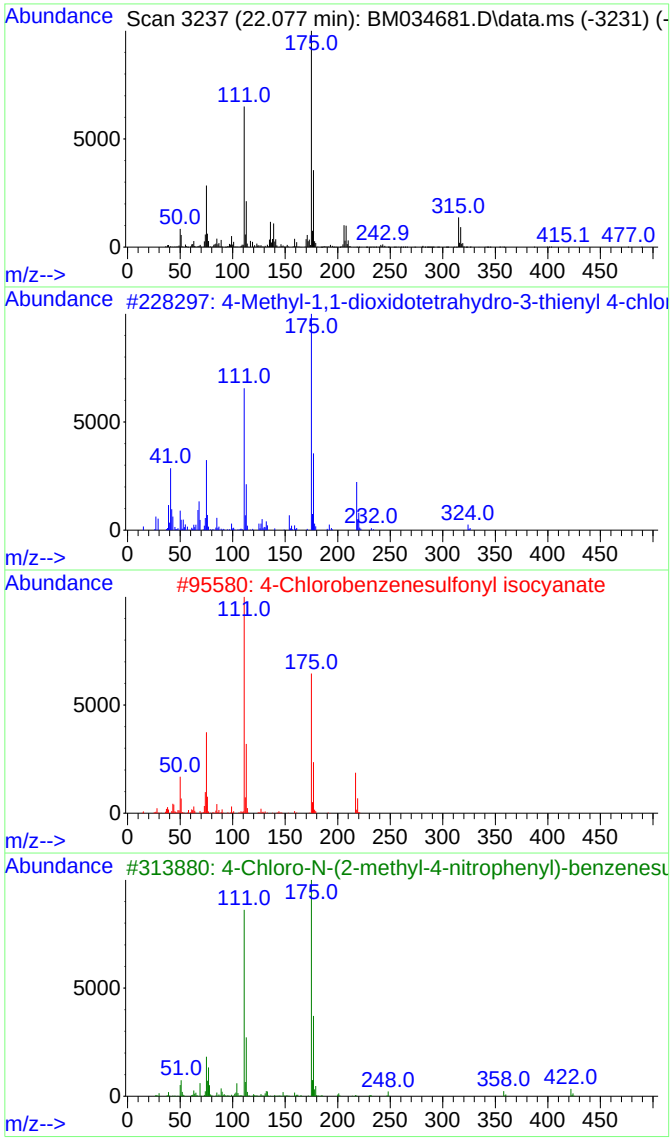
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

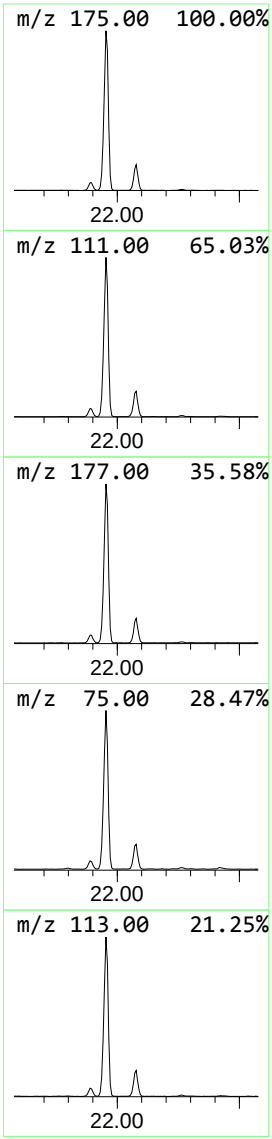
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 4-Chlorobenzenesulfonyl iso... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.077	36.85 ng/ul	2483170	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-1,1-dioxidotetrahydro-3...	324	C11H13ClO5S2	1000332-40-7	64
2	4-Chlorobenzenesulfonyl isocyanate	217	C7H4ClNO3S	005769-15-3	59
3	4-Chloro-N-(2-methyl-4-nitrophen...	422	C15H10ClF3N2O5S	1000374-78-8	45
4	Benzenesulfonyl chloride, 4-chloro-	210	C6H4Cl2O2S	000098-60-2	42
5	Benzene, 1-chloro-4-[(chlorometh...	224	C7H6Cl2O2S	005943-04-4	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

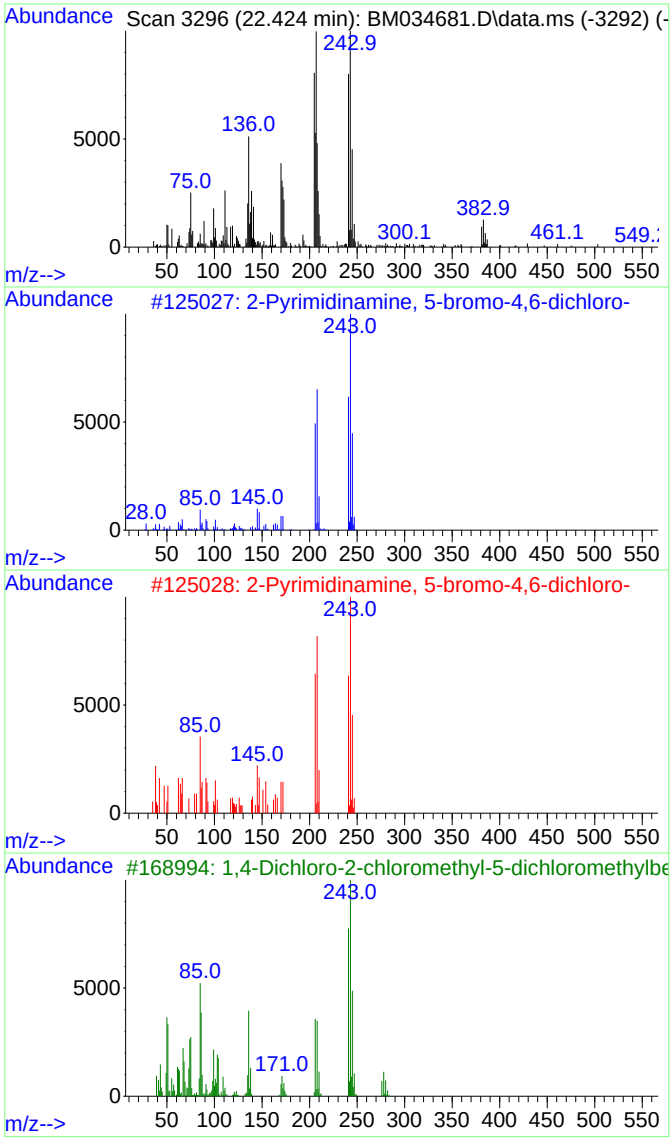
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

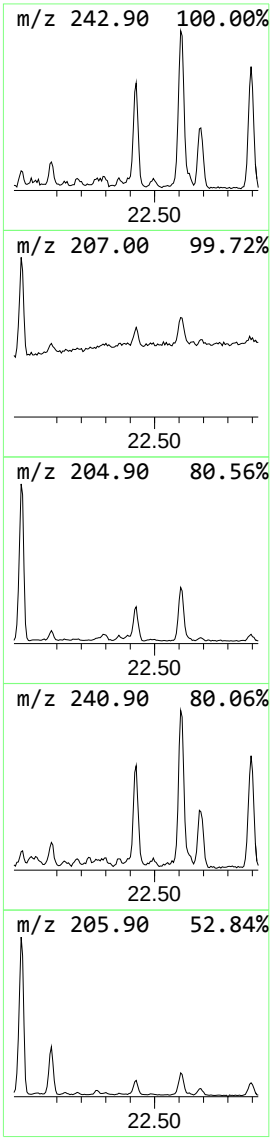
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 2-Pyrimidinamine, 5-bromo-4... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.424	5.78 ng/ul	389771	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pyrimidinamine, 5-bromo-4,6-di...	241	C4H2BrCl2N3	007781-26-2	60
2	2-Pyrimidinamine, 5-bromo-4,6-di...	241	C4H2BrCl2N3	007781-26-2	50
3	1,4-Dichloro-2-chloromethyl-5-di...	276	C8H5Cl5	1000306-27-6	47
4	Phthalide, 4,5,6,7-tetrachloro-	270	C8H2Cl4O2	027355-22-2	22
5	2-Bromo-4-chloroaniline	205	C6H5BrClN	000873-38-1	18





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

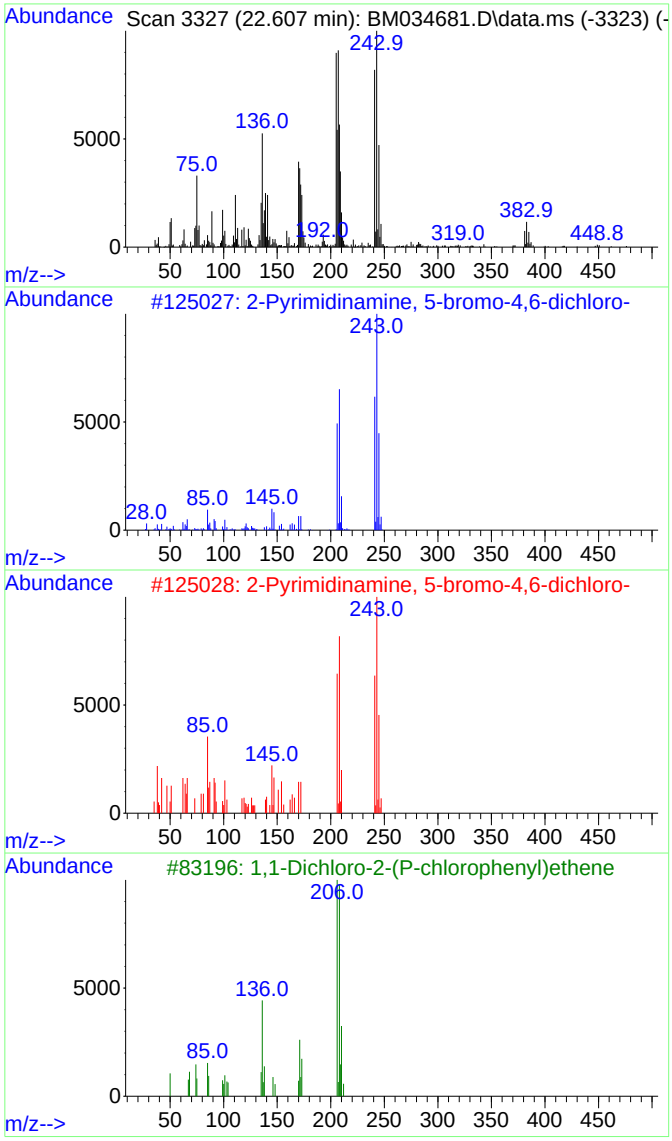
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

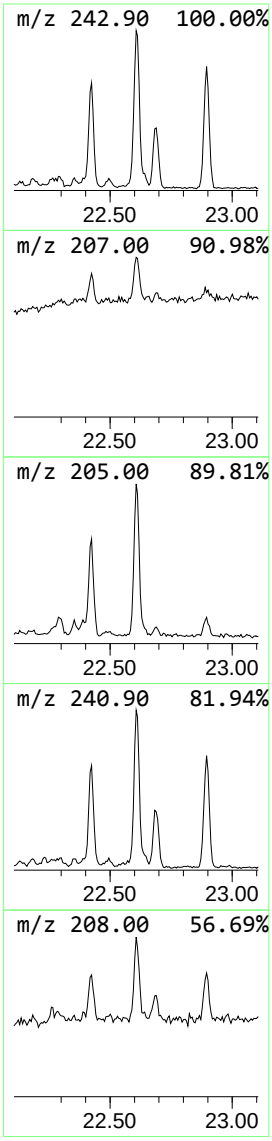
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 1,1-Dichloro-2-(P-chlorophe... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.607	9.27 ng/ul	624350	Chrysene-d12	21.454

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pyrimidinamine, 5-bromo-4,6-di...	241	C4H2BrCl2N3	007781-26-2	64
2	2-Pyrimidinamine, 5-bromo-4,6-di...	241	C4H2BrCl2N3	007781-26-2	64
3	1,1-Dichloro-2-(P-chlorophenyl)e...	206	C8H5Cl3	005263-17-2	44
4	Phthalide, 4,5,6,7-tetrachloro-	270	C8H2Cl4O2	027355-22-2	38
5	4-Trichloromethylphenyl phenylsu...	382	C14H10Cl4O2S	1000254-14-4	35





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Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

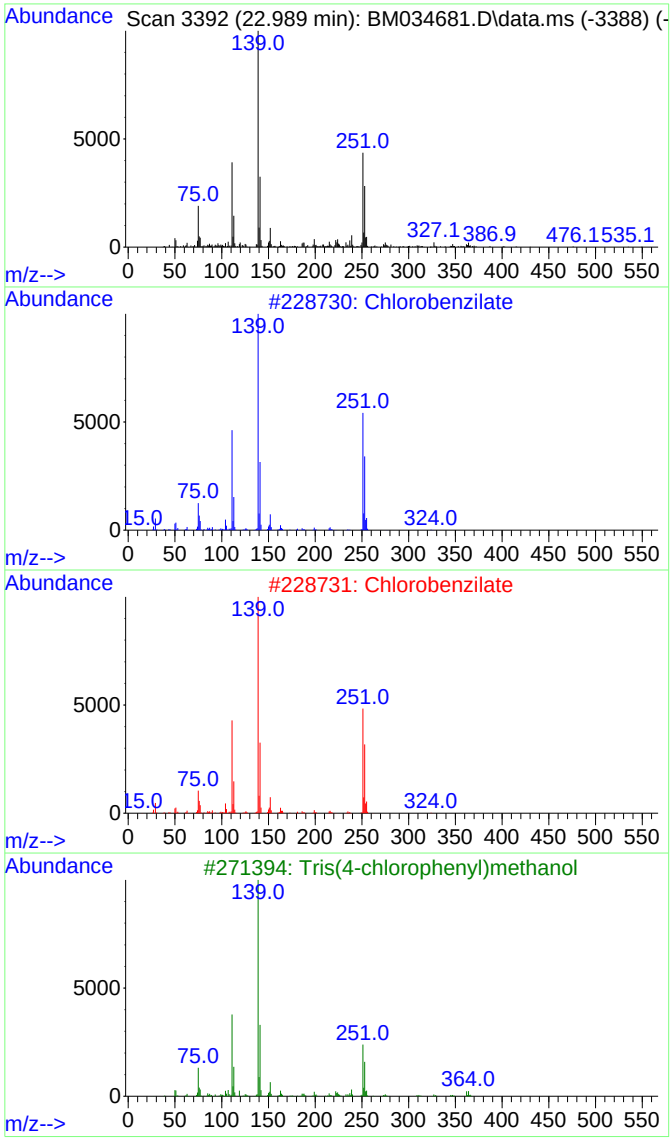
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

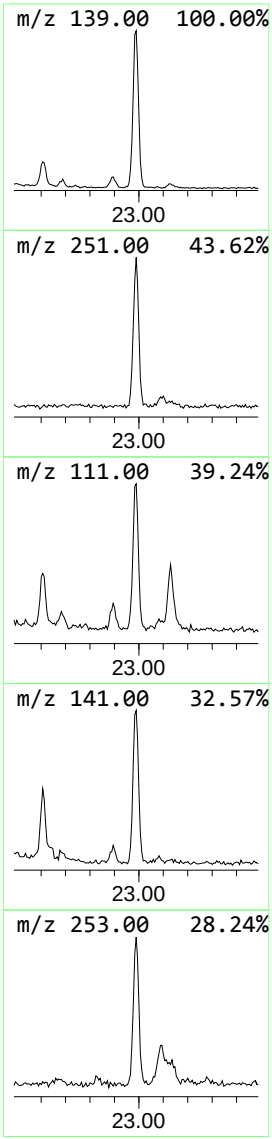
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Chlorobenzilate Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.989	4.21 ng/ul	371474	Perylene-d12	23.806

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chlorobenzilate	324	C16H14Cl2O3	000510-15-6	95
2	Chlorobenzilate	324	C16H14Cl2O3	000510-15-6	95
3	Tris(4-chlorophenyl)methanol	362	C19H13Cl3O	003010-80-8	93
4	4-Chloro-N-(3-nitrophenyl)benzamide	276	C13H9ClN2O3	006282-15-1	91
5	Chloropropylate	338	C17H16Cl2O3	005836-10-2	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034681.D
Acq On : 14 Apr 2022 23:26
Operator : CG/JU
Sample : N2316-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.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
Benzene, 1,4-di...	13.166	18.6	ng/ul	1345890	3	14.513	1451110	20.0
1,2,2-Trichloro...	14.672	69.7	ng/ul	5060030	3	14.513	1451110	20.0
unknown-01	15.683	7.1	ng/ul	515388	3	14.513	1451110	20.0
Benzene, 1-chlo...	15.736	11.3	ng/ul	819040	3	14.513	1451110	20.0
unknown-02	16.172	13.8	ng/ul	1063010	4	17.254	1537460	20.0
3,5-Dichloroben...	16.283	68.5	ng/ul	5267420	4	17.254	1537460	20.0
o,p'-DDE	19.401	24.7	ng/ul	1664210	5	21.454	1347680	20.0
Mitotane	19.554	25.0	ng/ul	1684590	5	21.454	1347680	20.0
Hydroxyzine	19.583	13.6	ng/ul	913708	5	21.454	1347680	20.0
unknown-03	19.724	22.3	ng/ul	1500860	5	21.454	1347680	20.0
p,p'-DDE	19.789	180.4	ng/ul	12153100	5	21.454	1347680	20.0
m,p'-DDD	19.895	176.7	ng/ul	11905500	5	21.454	1347680	20.0
unknown-04	19.936	34.1	ng/ul	2294790	5	21.454	1347680	20.0
5-[2-(5-Oxopent...	20.089	32.7	ng/ul	2203910	5	21.454	1347680	20.0
9H-Carbazole, 3...	20.124	106.9	ng/ul	7203320	5	21.454	1347680	20.0
Benzene, 1,1'-s...	20.260	10.1	ng/ul	679339	5	21.454	1347680	20.0
o,p'-DDT	20.366	3022.4	ng/ul	203659000	5	21.454	1347680	20.0
p,p'-DDT	20.530	406.4	ng/ul	27387000	5	21.454	1347680	20.0
unknown-05	20.760	5433.9	ng/ul	366156000	5	21.454	1347680	20.0
1,1-Dichloro-2,...	20.901	188.7	ng/ul	12715300	5	21.454	1347680	20.0
3,5-Dibromo-2-c...	21.160	12.3	ng/ul	828059	5	21.454	1347680	20.0
unknown-06	21.307	5.8	ng/ul	391892	5	21.454	1347680	20.0
4-Methyl-1,1-di...	21.889	10.8	ng/ul	726789	5	21.454	1347680	20.0
4-Chloro-N-(2-m...	21.954	197.0	ng/ul	13276000	5	21.454	1347680	20.0
4-Chlorobenzene...	22.077	36.9	ng/ul	2483170	5	21.454	1347680	20.0
2-Pyrimidinamin...	22.424	5.8	ng/ul	389771	5	21.454	1347680	20.0
1,1-Dichloro-2-...	22.607	9.3	ng/ul	624350	5	21.454	1347680	20.0
Chlorobenzilate	22.989	4.2	ng/ul	371474	6	23.806	1765750	20.0