LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : GBNJ4ME

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-BM120122.M
Title : SVOA CALIBRATION

Signal : TIC: BM037859.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.245	649	656	666	rBV	782282	1230895	0.57%	0.379%
2	7.410	678	684	692	rBV	659988	1034552	0.48%	0.318%
3	7.616	712	719	727	rBV	790193	1239018	0.57%	0.381%
4	8.092	793	800	807	rBV	904938	1405321	0.65%	0.433%
5	8.798	911	920	932	rBV	946700	1535759	0.71%	0.473%
6	9.263	992	999	1006	rBV	801489	1276806	0.59%	0.393%
7	9.986	1114	1122	1132	rBV	632921	1036348	0.48%	0.319%
8	10.528	1207	1214	1225	rBV	927172	1520292	0.70%	0.468%
9	10.910	1269	1279	1288	rBV	1242105	2075534	0.96%	0.639%
10	11.051	1295	1303	1317	rBV	488996	905283	0.42%	0.279%
11	14.127	1817	1826	1834	rBV	1864827	2478514	1.14%	0.763%
12	14.416	1868	1875	1902	rBV	1559788	3039922	1.40%	0.936%
13	14.721	1918	1927	1952	rBV	1504713	2721181	1.26%	0.838%
14	14.910	1952	1959	1976	rBV	673003	1056500	0.49%	0.325%
15	15.351	2022	2034	2054	rVV	16196293	29532187	13.64%	9.092%
16	15.710	2088	2095	2106	rVB	2687940	3865752	1.79%	1.190%
17	15.821	2108	2114	2128	rVV	680672	1060099	0.49%	0.326%
18	17.210	2324	2350	2354	rBV	42911145	216495739	100.00%	66.650%
19	17.251	2354	2357	2386	rVB	718006	2116469	0.98%	0.652%
20	17.480	2391	2396	2403	rVV	2175716	2988666	1.38%	0.920%
21	17.574	2406	2412	2418	rVV	2704891	3652891	1.69%	1.125%
22	18.015	2482	2487	2491	rBV	170095	240532	0.11%	0.074%
23	18.068	2491	2496	2502	rVB	836391	1093456	0.51%	0.337%
24	18.333	2536	2541	2551	rBV	654670	964604	0.45%	0.297%
25	19.598	2753	2756	2763	rBV	170016	262655	0.12%	0.081%
26	19.839	2791	2797	2805	rBV	2804220	3692188	1.71%	1.137%
27	21.309	3043	3047	3054	rBV	827748	1060946	0.49%	0.327%
28	21.621	3096	3100	3108	rVB	1775772	2319769	1.07%	0.714%
29	22.703	3280	3284	3289	rBV	419444	664351	0.31%	0.205%
30	22.892	3313	3316	3325	rVB2	324748	681625	0.31%	0.210%
31	23.950	3490	3496	3502	rBV2	1264110	2492261	1.15%	0.767%
32	24.050	3507	3513	3518	rVV	3254232	6665356	3.08%	2.052%
33	24.109	3519	3523	3530	rVB2	1265170	2463670	1.14%	0.758%
34	24.191	3532	3537	3542	rVV	1067930	1967605	0.91%	0.606%
35	24.250	3543	3547	3557	rVB3	577168	1316690	0.61%	0.405%
36	24.444	3575	3580	3601	rVB4	566055	1595018	0.74%	0.491%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

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-BM120122.M
Title : SVOA CALIBRATION

37	25.650	3779	3785	3789	rBV3	1030314	2569771	1.19%	0.791%
38	25.709	3790	3795	3811	rVB2	2010738	5151283	2.38%	1.586%
39	25.891	3818	3826	3834	rBV	1257869	3130263	1.45%	0.964%
40	26.021	3842	3848	3867	rVB5	1505778	4224864	1.95%	1.301%

Sum of corrected areas: 324824635

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
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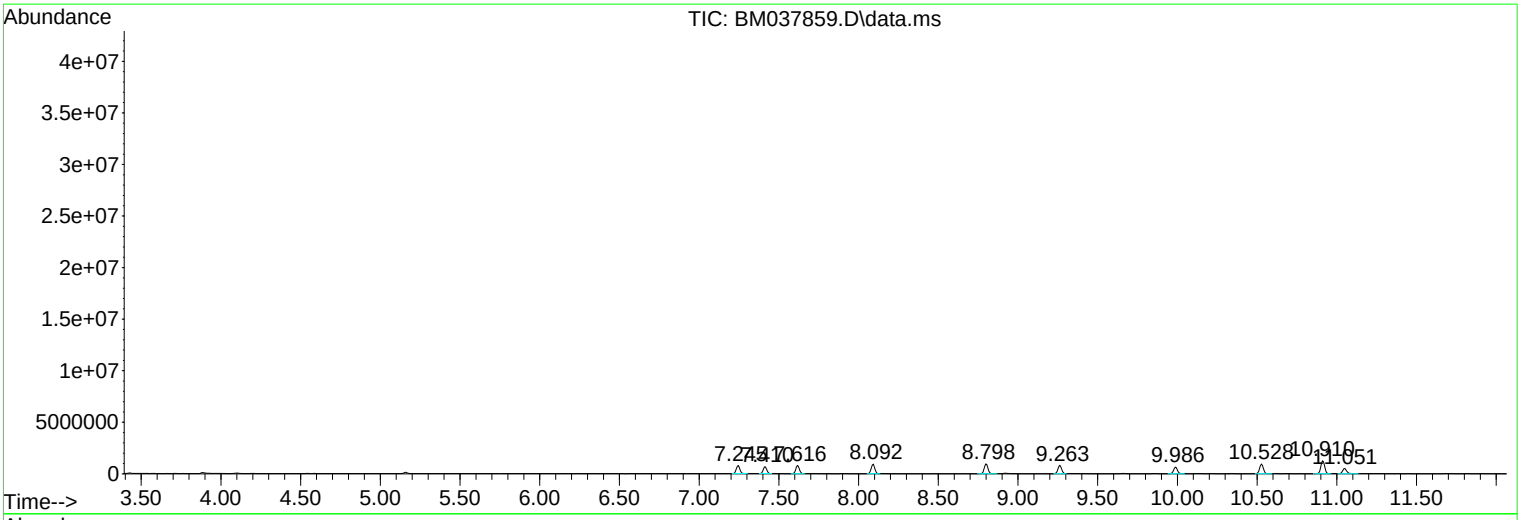
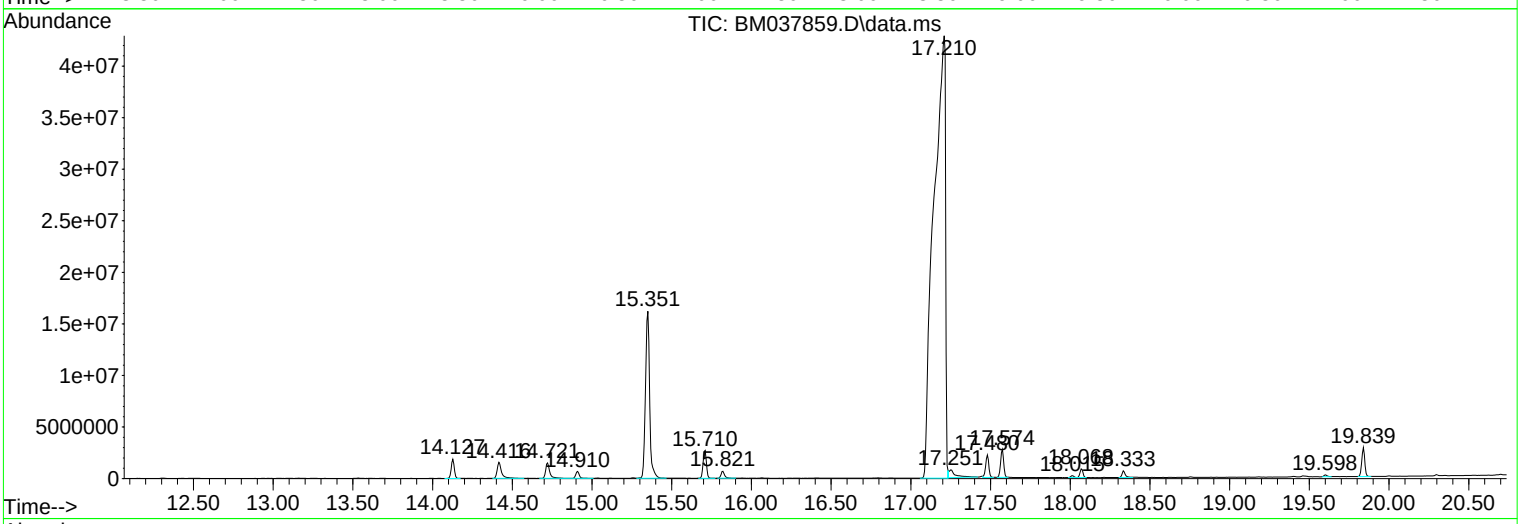
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

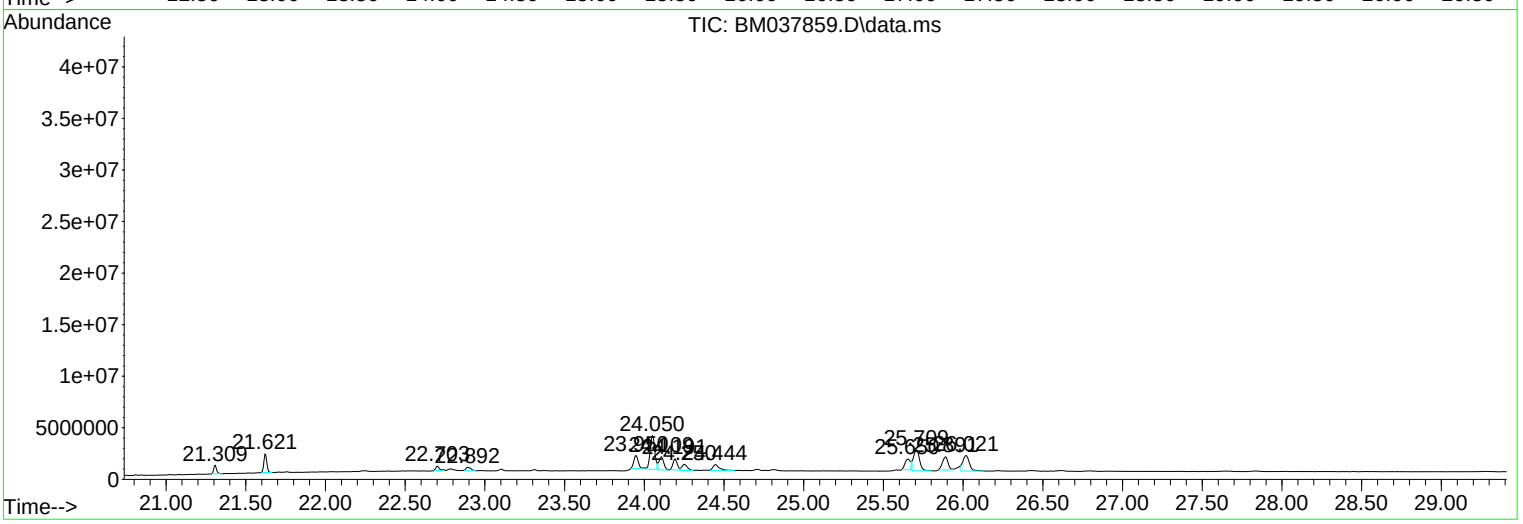
Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

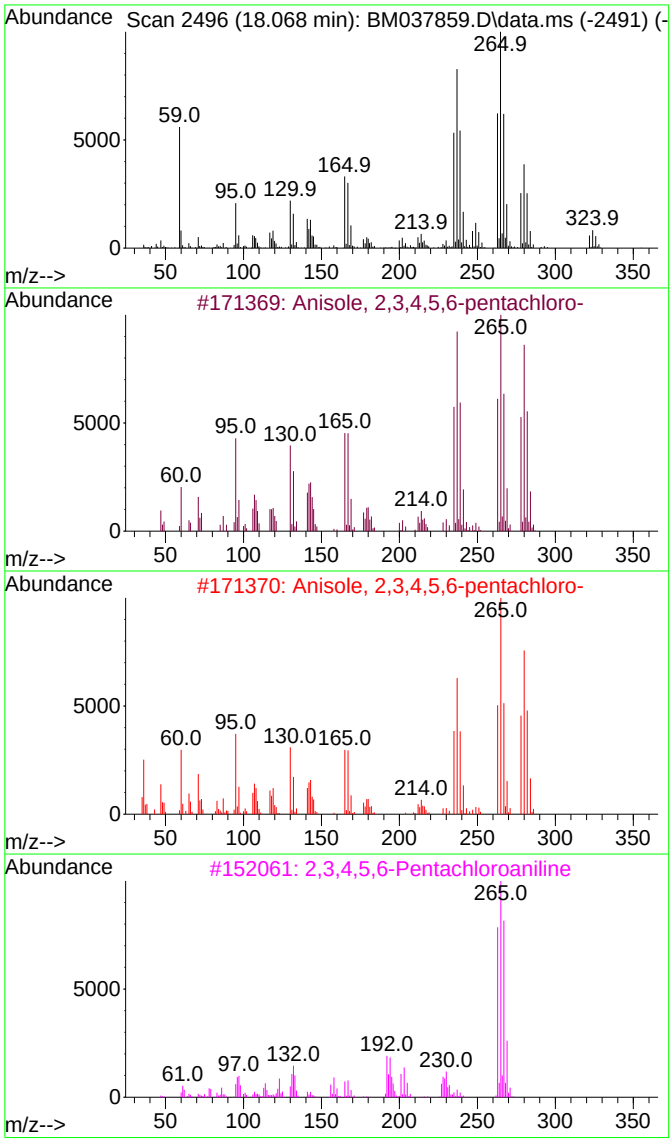
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

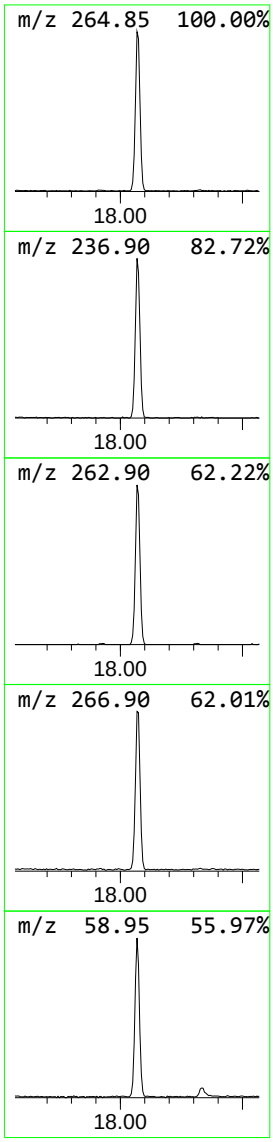
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Anisole, 2,3,4,5,6-pentachl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.068	7.32 ng/ul	1093460	Phenanthrene-d10	17.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	95
2	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	78
3	2,3,4,5,6-Pentachloroaniline	263	C6H2Cl5N	000527-20-8	38
4	2,3,4,5,6-Pentachloroaniline	263	C6H2Cl5N	000527-20-8	30
5	Rhodium, acetylanilinato-bis(eth...	293	C12H16NORh	1000153-76-6	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

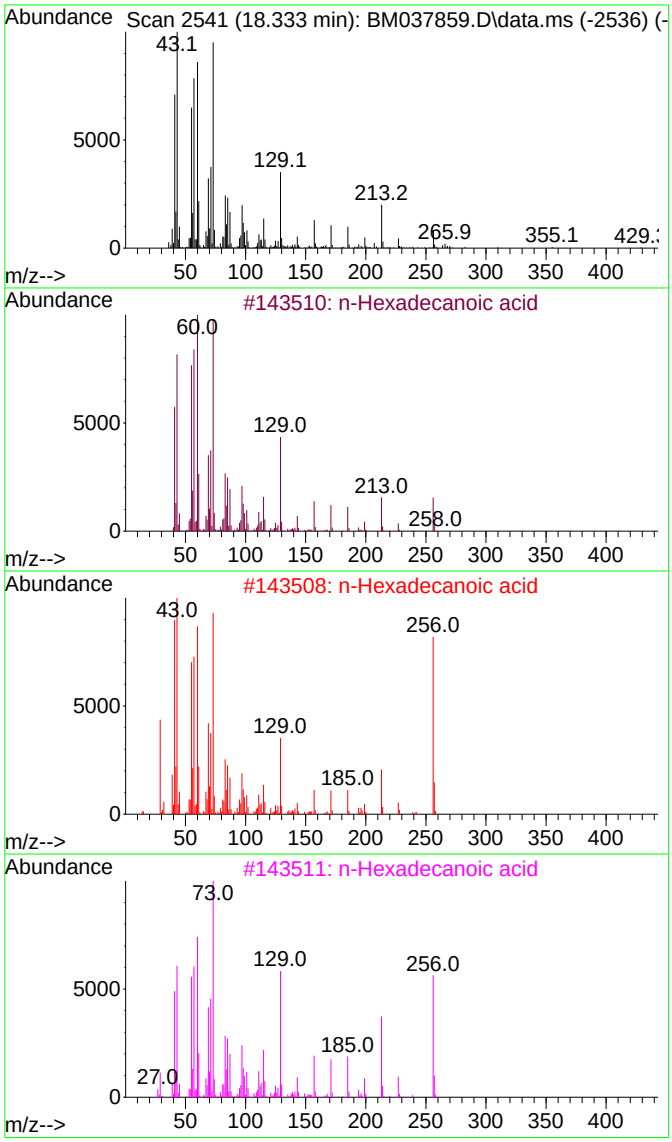
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

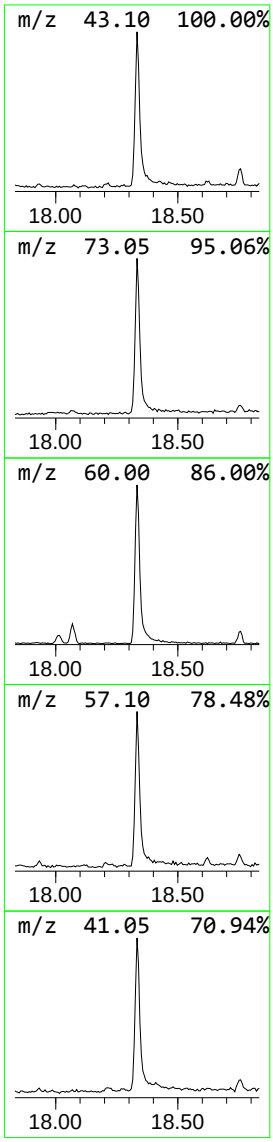
****

Peak Number 2 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.333	6.46 ng/ul	964604	Phenanthrene-d10	17.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5	Tridecanoic acid	214	C13H26O2	000638-53-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

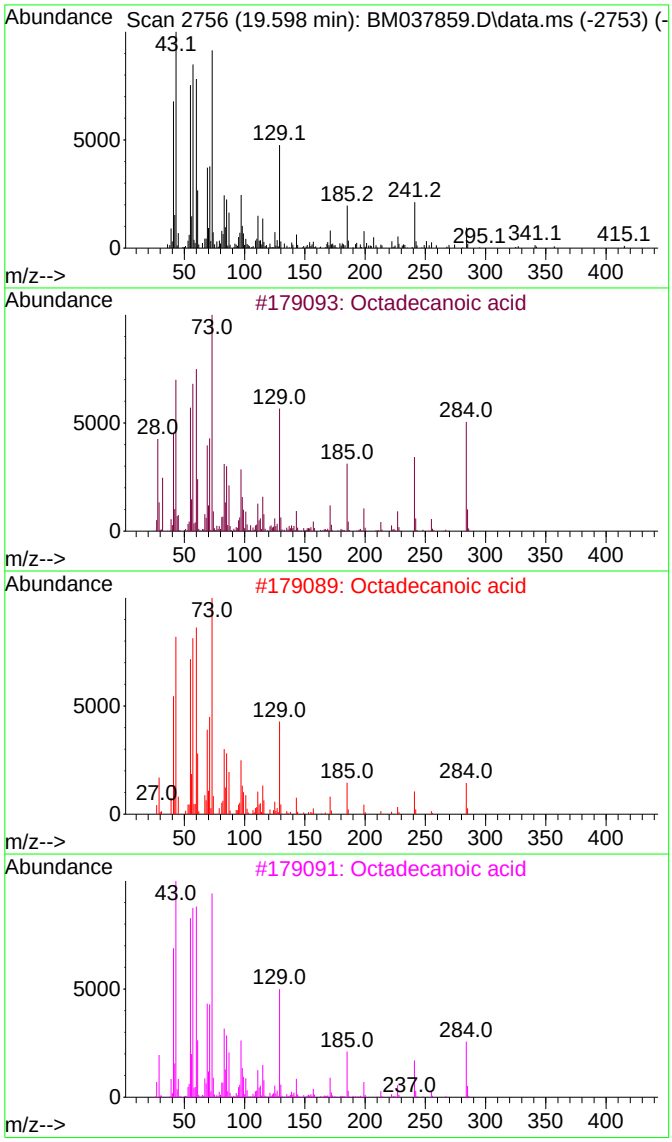
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

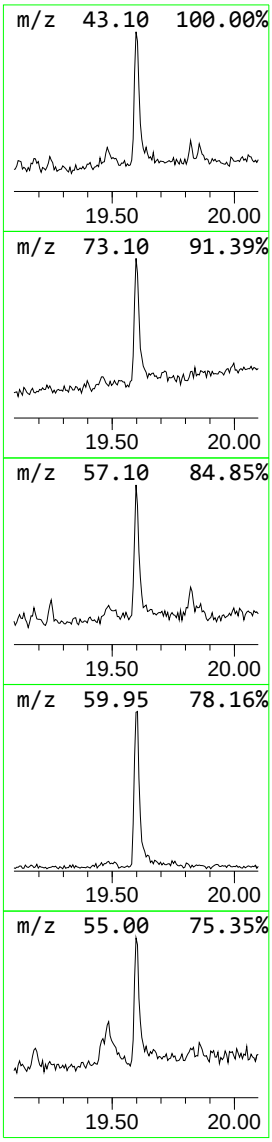
****

Peak Number 3 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.598	2.26 ng/ul	262655	Chrysene-d12	21.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	96
3	Octadecanoic acid	284	C18H36O2	000057-11-4	93
4	Octadecanoic acid	284	C18H36O2	000057-11-4	90
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

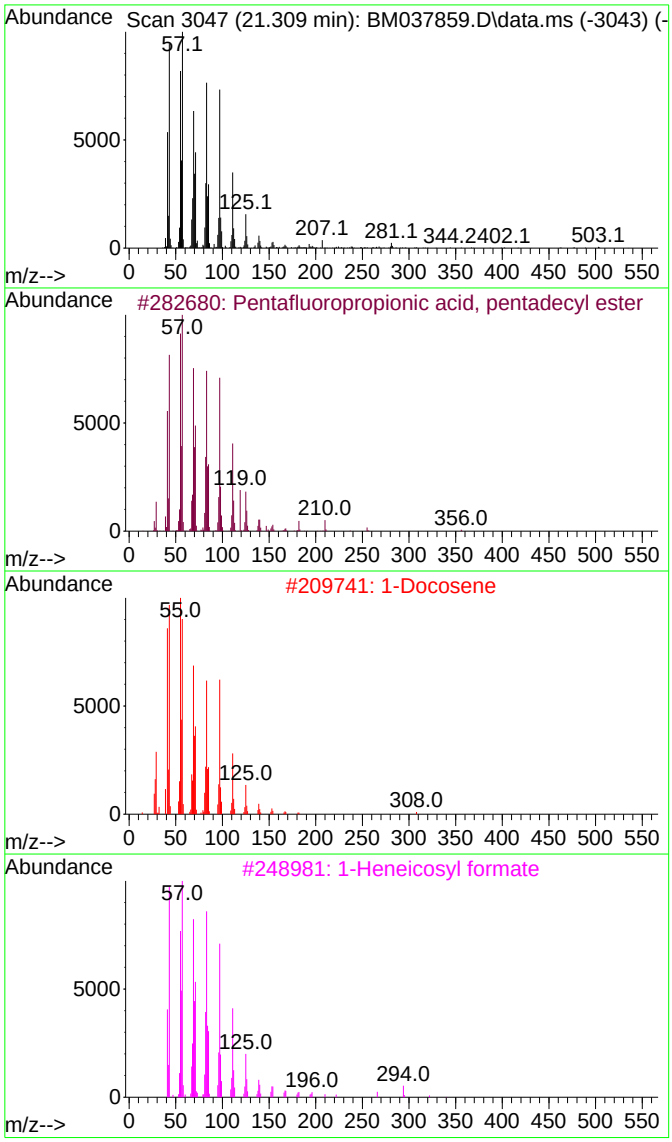
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

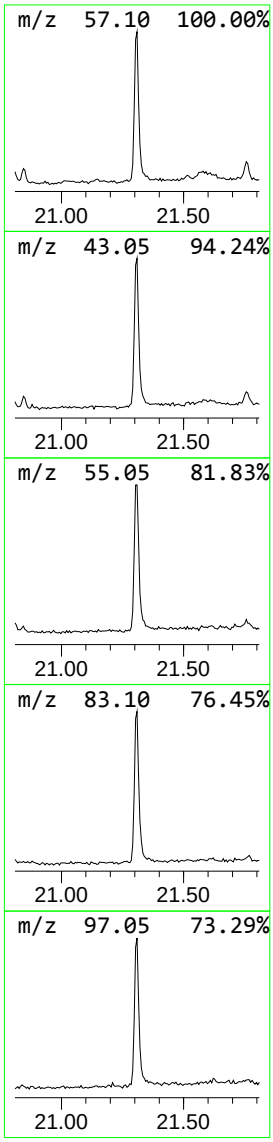
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentafluoropropionic acid, ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.309	9.15 ng/ul	1060950	Chrysene-d12	21.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
2	1-Docosene	308	C22H44	001599-67-3	94
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	1-Tetracosene	336	C24H48	010192-32-2	94
5	1-Hexacosene	364	C26H52	018835-33-1	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

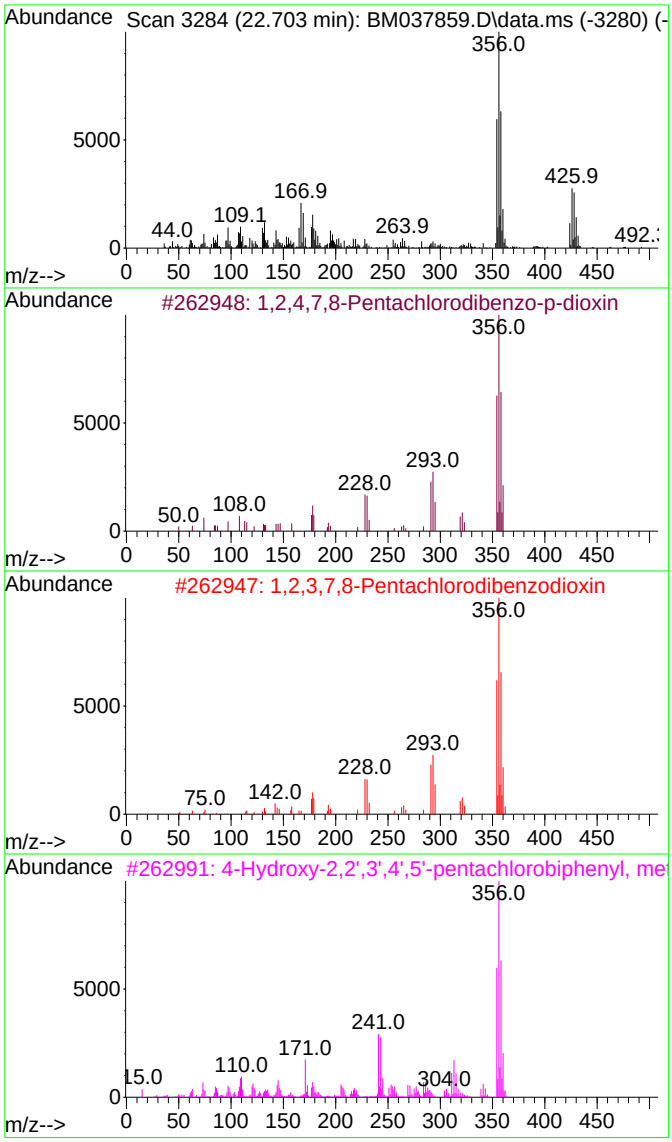
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

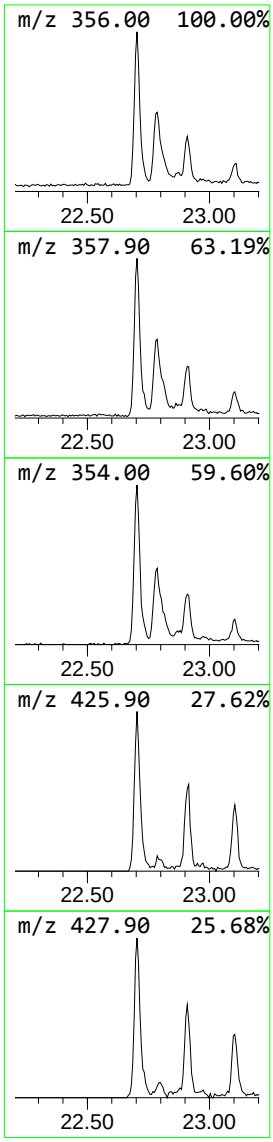
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,2,4,7,8-Pentachlorodibenz... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.703	5.73 ng/ul	664351	Chrysene-d12	21.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,7,8-Pentachlorodibenzo-p-d...	354	C12H3Cl5O2	058802-08-7	91
2	1,2,3,7,8-Pentachlorodibenzodioxin	354	C12H3Cl5O2	040321-76-4	83
3	4-Hydroxy-2,2',3',4',5'-pentachl...	354	C13H7Cl5O	1000447-88-3	64
4	4-Hydroxy-2,2',3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-4	64
5	4-Hydroxy-2',3,3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-6	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

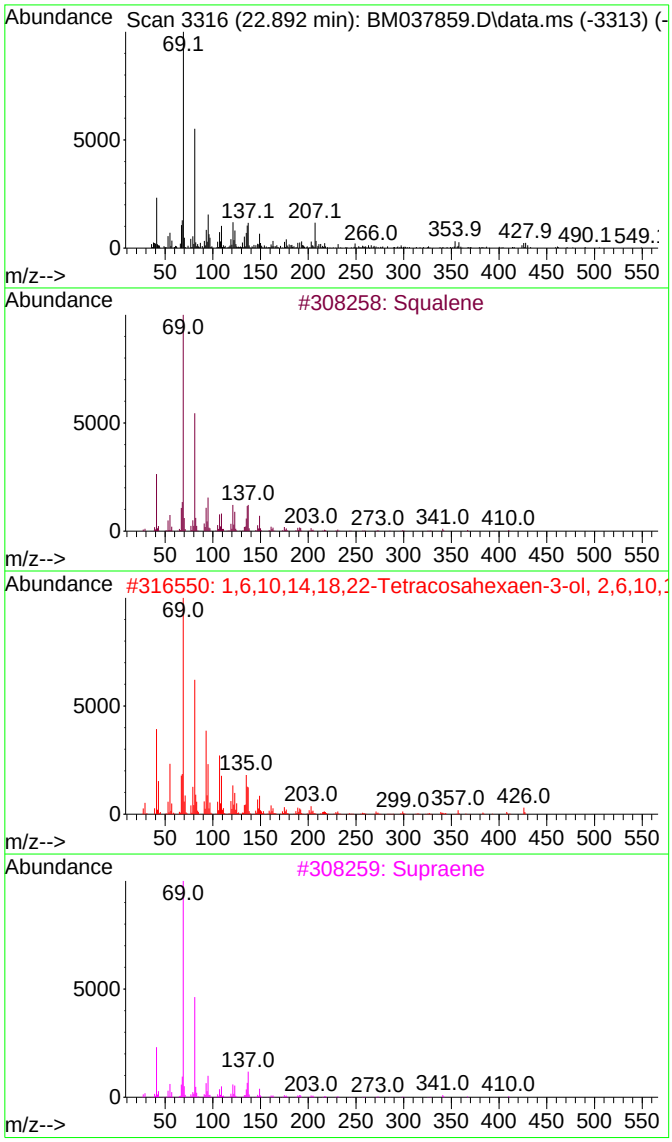
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

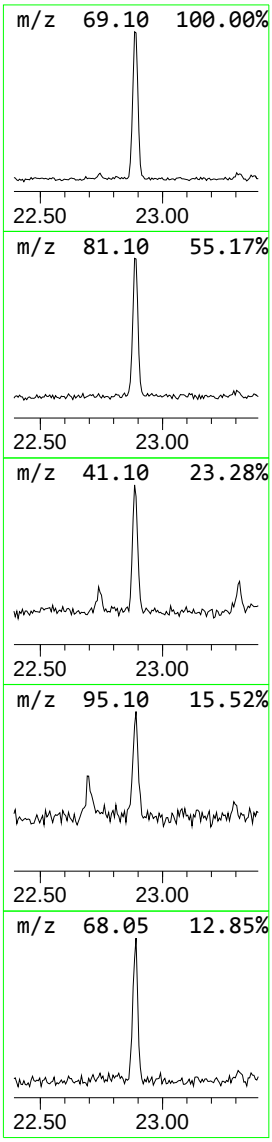
****

Peak Number 6 Squalene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.892	5.53 ng/ul	681625	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	83
2	1,6,10,14,18,22-Tetracosahexaen-...	426	C30H50O	054159-46-5	81
3	Supraene	410	C30H50	007683-64-9	80
4	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	72
5	Squalene	410	C30H50	000111-02-4	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

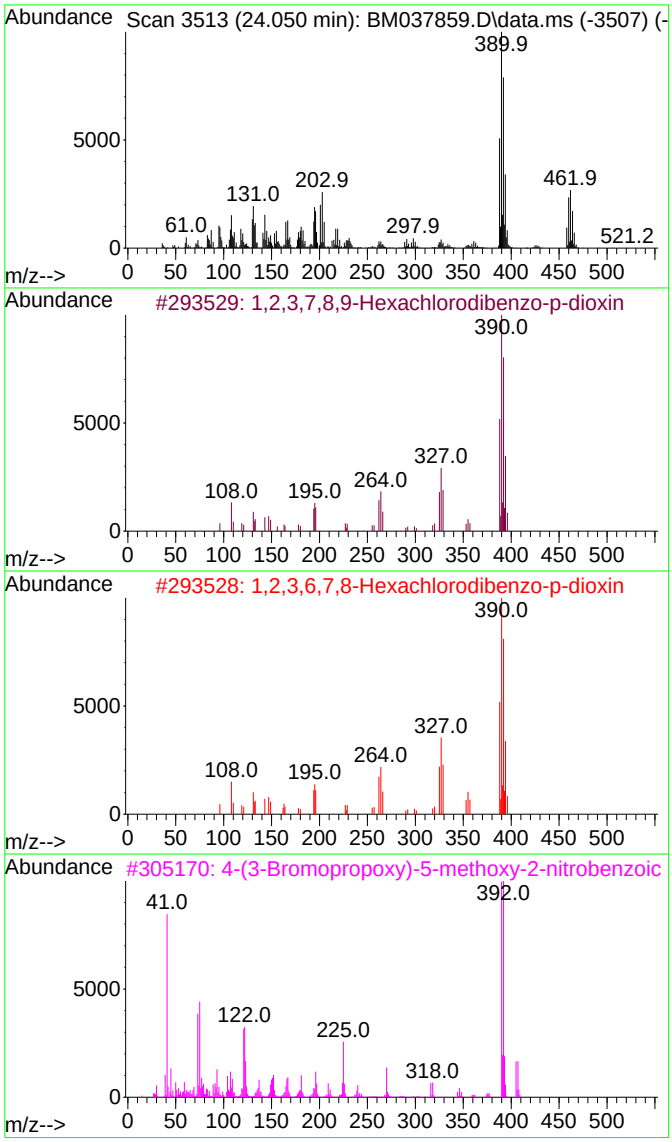
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

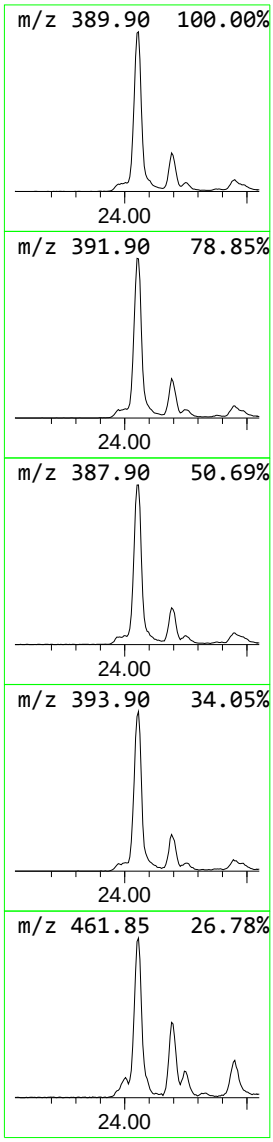
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,2,3,7,8,9-Hexachlorodiben... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.050	54.11 ng/ul	6665360	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,7,8,9-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	019408-74-3	89
2	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	89
3	4-(3-Bromopropoxy)-5-methoxy-2-n...	405	C14H20BrNO6Si	1000471-26-2	37
4	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	32
5	4H-1-Benzopyran-2-carboxylic aci...	390	C12H8Br2O5	030113-88-3	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

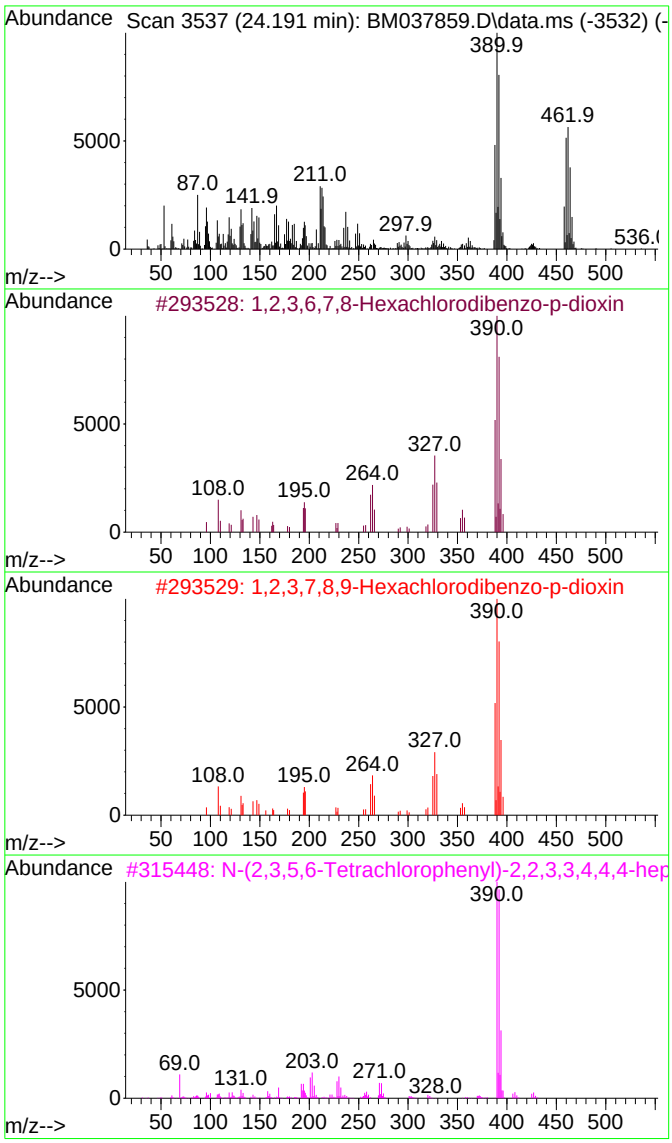
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

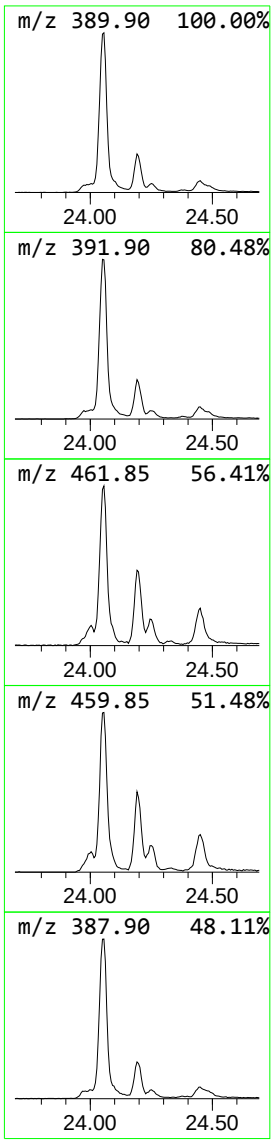
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,2,3,6,7,8-Hexachlorodiben... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.191	15.97 ng/ul	1967610	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	78
2	1,2,3,7,8,9-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	019408-74-3	60
3	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	38
4	Thieno[2,3-b]thiophene, 2-bromo-...	390	C11H8Br2N2S2	1000268-03-5	25
5	Pregn-20-yn-6-one, 17-(acetyloxy...	390	C23H31FO4	057194-88-4	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

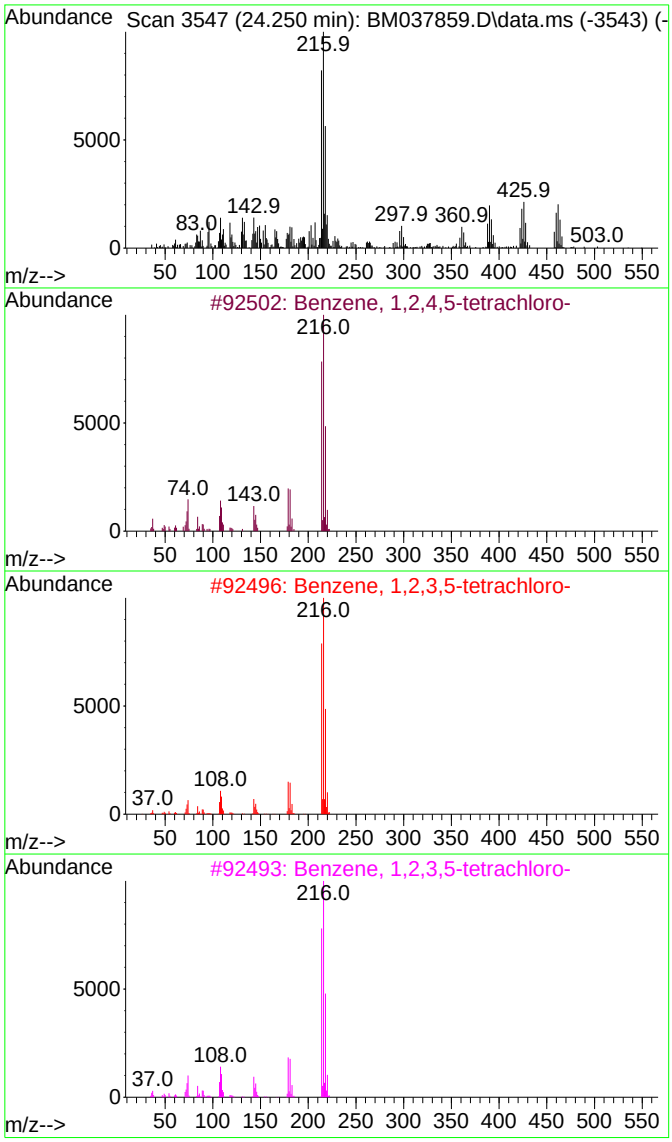
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

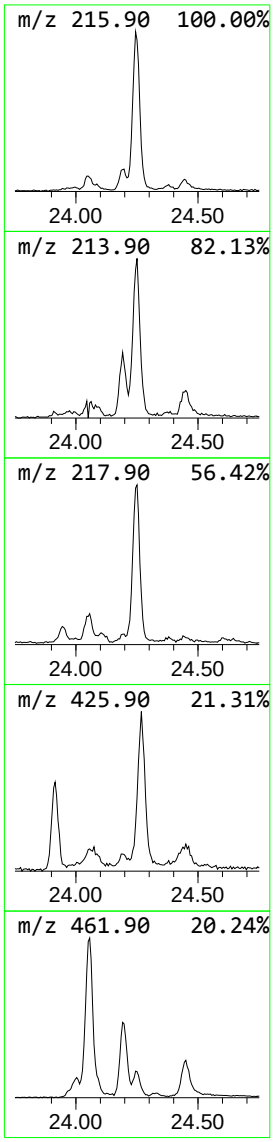
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,4,5-tetrachloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.250	10.69 ng/ul	1316690	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	78
2	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	78
3	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	78
4	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	78
5	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	60





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Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

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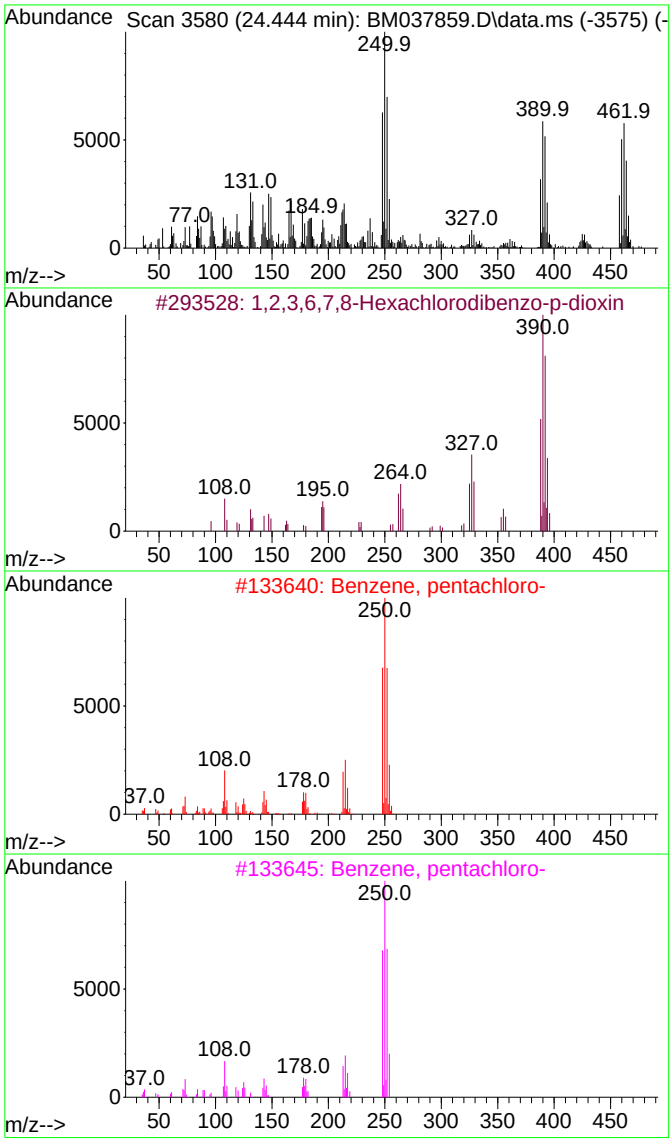
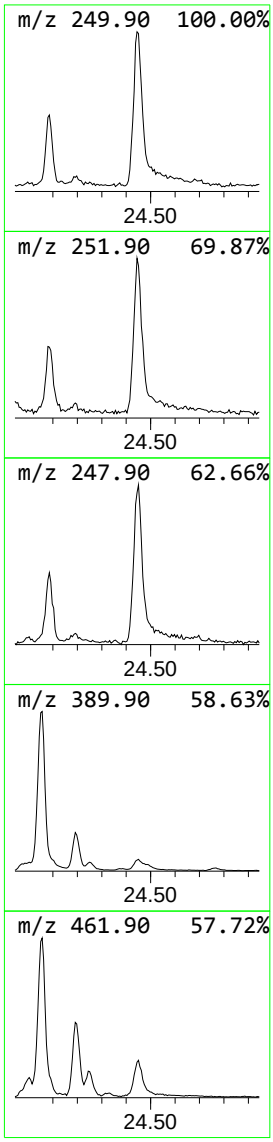
Peak Number 10 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.444	12.95 ng/ul	1595020	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	56
2	Benzene, pentachloro-	248	C6HCl5	000608-93-5	38
3	Benzene, pentachloro-	248	C6HCl5	000608-93-5	38
4	Benzene, pentachloro-	248	C6HCl5	000608-93-5	38
5	Benzene, pentachloro-	248	C6HCl5	000608-93-5	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

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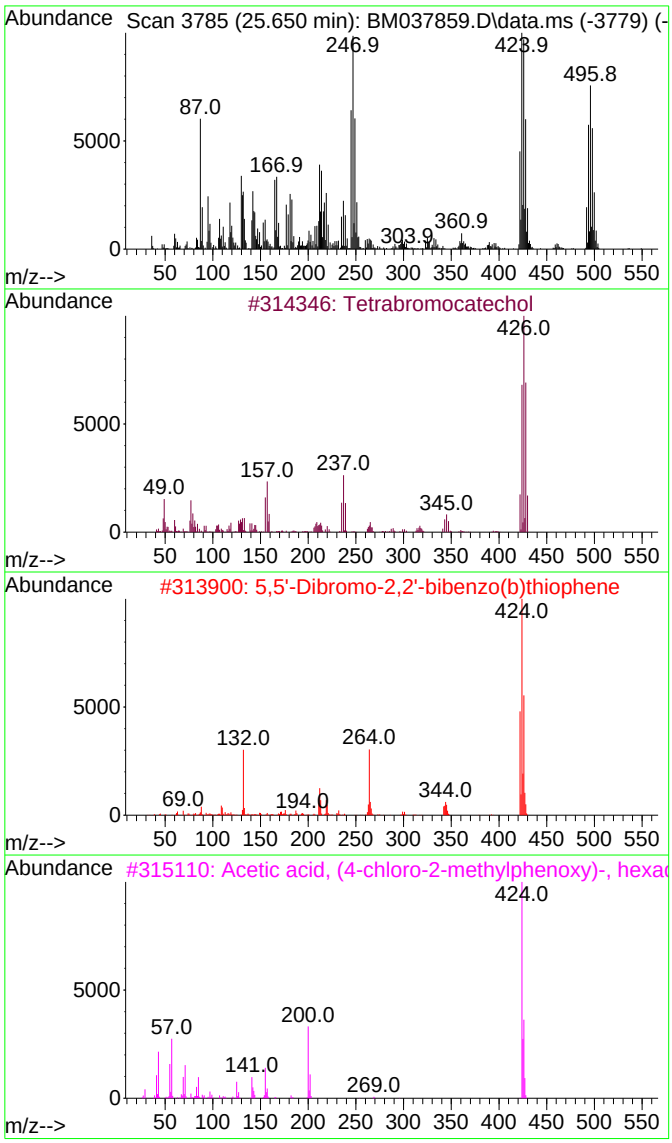
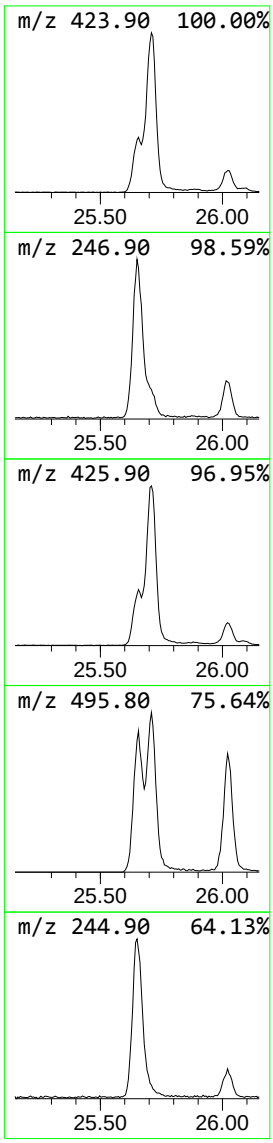
Peak Number 11 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.650	20.86 ng/ul	2569770	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrabromocatechol	422	C6H2Br4O2	000488-47-1	35
2	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	20
3	Acetic acid, (4-chloro-2-methylp...	424	C25H41ClO3	1000415-16-9	11
4	Dibenzo(b,f)(1,5)diazocine, 2,8-...	426	C26H16Cl2N2	003646-61-5	10
5	4-(2-Bromophenyl)-2-[(2-bromophe...	424	C16H10Br2S2	1000444-65-0	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

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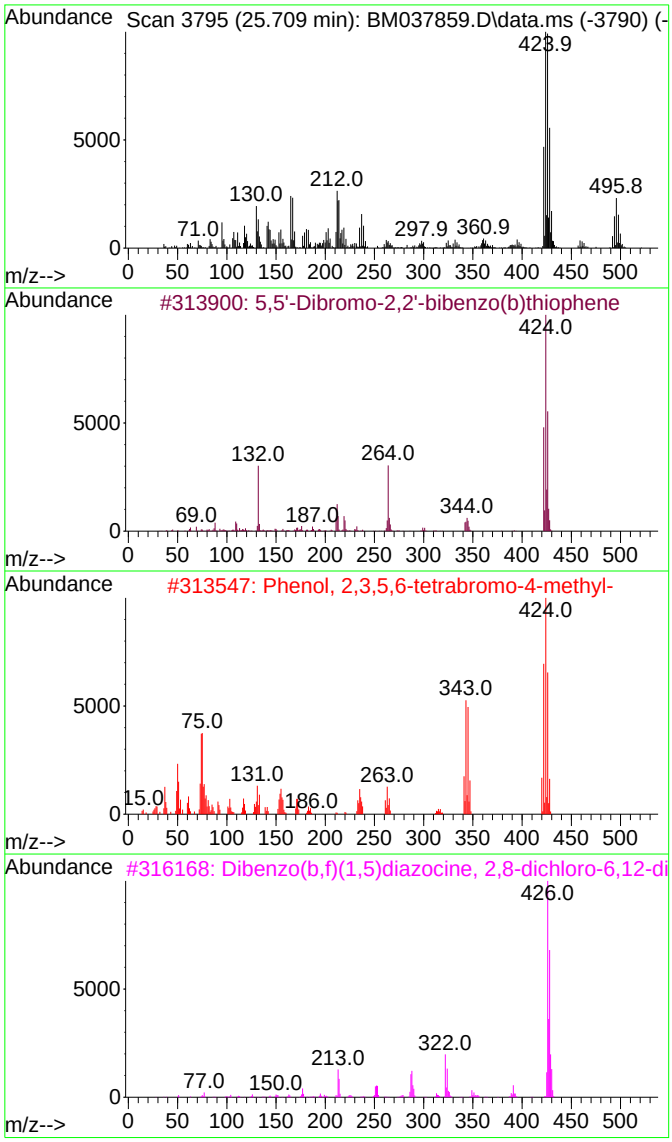
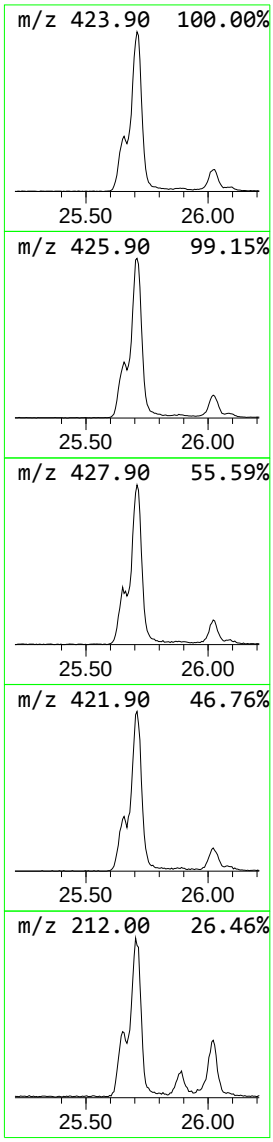
Peak Number 12 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.709	41.82 ng/ul	5151280	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	47
2	Phenol, 2,3,5,6-tetrabromo-4-met...	420	C7H4Br4O	037721-75-8	43
3	Dibenzo(b,f)(1,5)diazocine, 2,8-...	426	C26H16Cl2N2	003646-61-5	38
4	Phenol, 2,3,4,5-tetrabromo-6-met...	420	C7H4Br4O	000576-55-6	38
5	Phenol, 2,3,4,5-tetrabromo-6-met...	420	C7H4Br4O	000576-55-6	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

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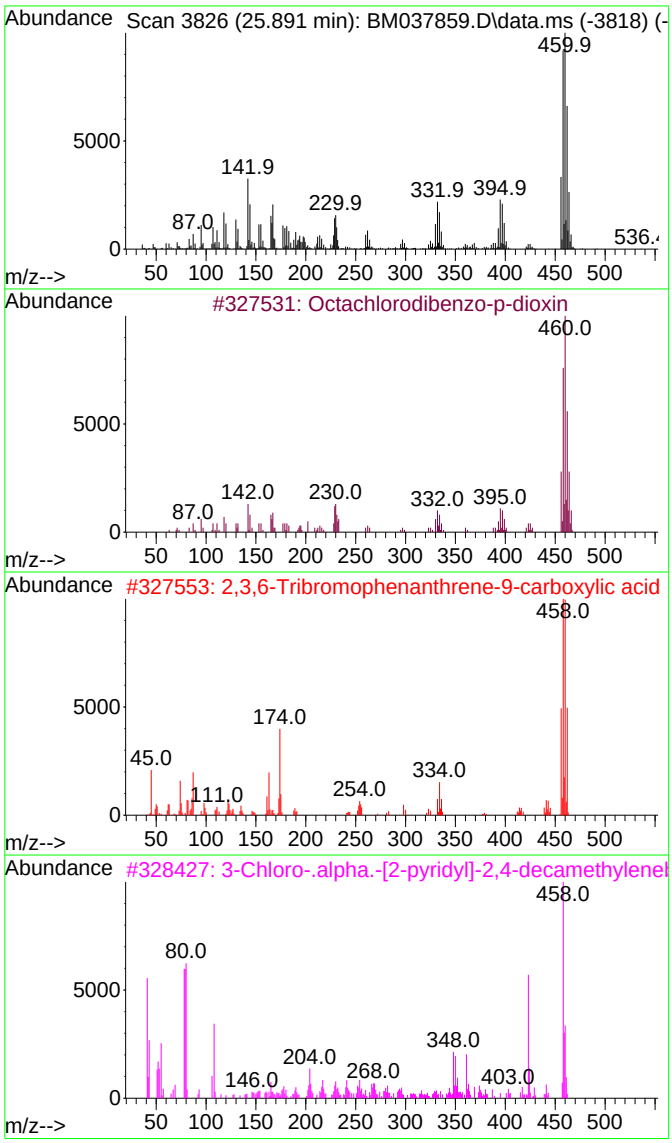
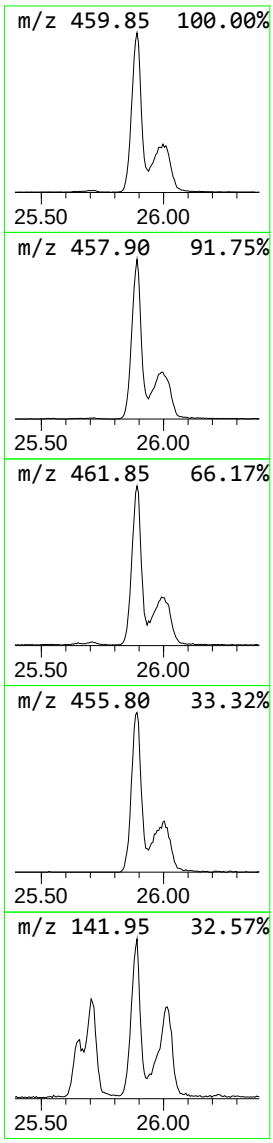
Peak Number 13 Octachlorodibenzo-p-dioxin Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.891	25.41 ng/ul	3130260	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	83
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	33
3	3-Chloro-.alpha.-[2-pyridyl]-2,4...	458	C29H31ClN2O	035158-92-0	14
4	1,1':4',1'':4'',1''':4''',1'''':...	458	C36H26	004499-83-6	9
5	Odoratin, 2TMS	458	C23H30O6Si2	1000502-41-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

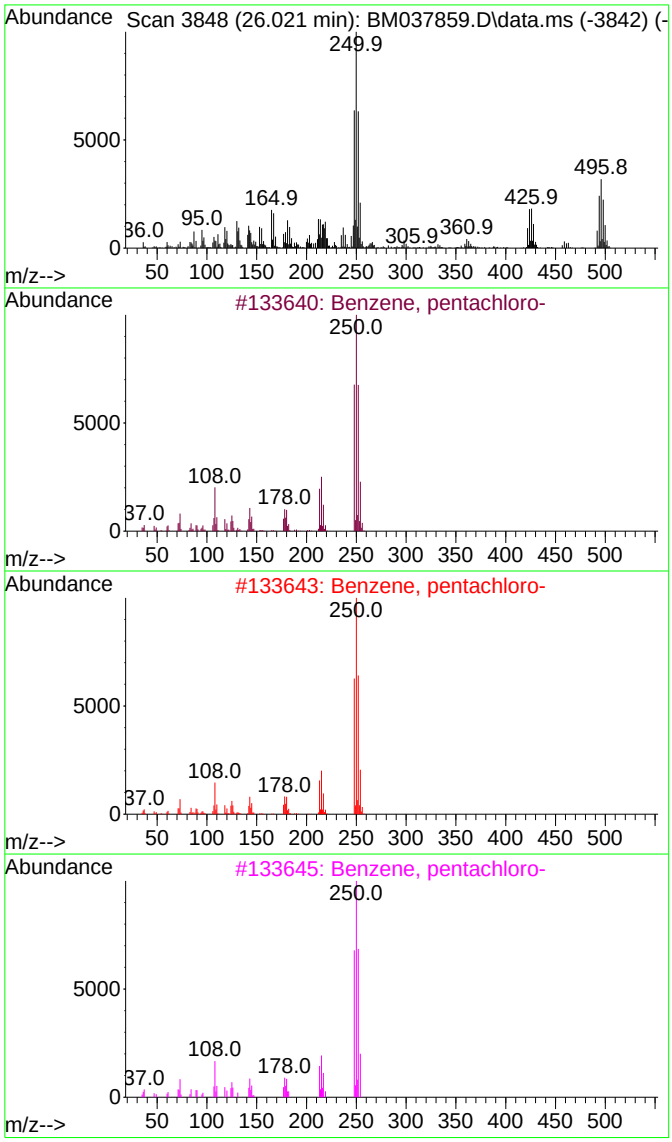
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

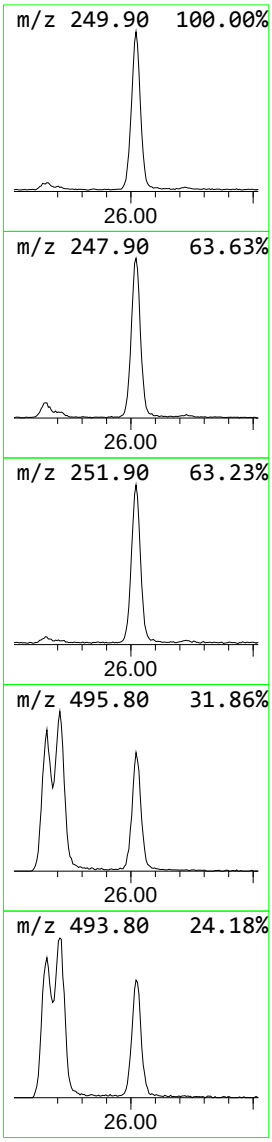
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, pentachloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.021	34.30 ng/ul	4224860	Perylene-d12	24.109

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentachloro-	248	C6HCl5	000608-93-5	78
2	Benzene, pentachloro-	248	C6HCl5	000608-93-5	78
3	Benzene, pentachloro-	248	C6HCl5	000608-93-5	60
4	4-Bromo-6-chlorobenzo[c][1,2,5]t...	248	C6H2BrClN2S	1215206-49-7	43
5	Fluvalinate	502	C26H22ClF3N2O3	069409-94-5	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037859.D
Acq On : 06 Dec 2022 18:07
Operator : CG/JU
Sample : N5738-15ME
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ4ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.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
Anisole, 2,3,4,...	18.068	7.3	ng/ul	1093460	4	17.480	2988670	20.0
n-Hexadecanoic ...	18.333	6.5	ng/ul	964604	4	17.480	2988670	20.0
Octadecanoic acid	19.598	2.3	ng/ul	262655	5	21.621	2319770	20.0
Pentafluoroprop...	21.309	9.2	ng/ul	1060950	5	21.621	2319770	20.0
1,2,4,7,8-Penta...	22.703	5.7	ng/ul	664351	5	21.621	2319770	20.0
Squalene	22.892	5.5	ng/ul	681625	6	24.109	2463670	20.0
1,2,3,7,8,9-Hex...	24.050	54.1	ng/ul	6665360	6	24.109	2463670	20.0
1,2,3,6,7,8-Hex...	24.191	16.0	ng/ul	1967610	6	24.109	2463670	20.0
Benzene, 1,2,4,...	24.250	10.7	ng/ul	1316690	6	24.109	2463670	20.0
unknown-01	24.444	12.9	ng/ul	1595020	6	24.109	2463670	20.0
unknown-02	25.650	20.9	ng/ul	2569770	6	24.109	2463670	20.0
unknown-03	25.709	41.8	ng/ul	5151280	6	24.109	2463670	20.0
Octachlorodiben...	25.891	25.4	ng/ul	3130260	6	24.109	2463670	20.0
Benzene, pentac...	26.021	34.3	ng/ul	4224860	6	24.109	2463670	20.0