LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

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-BP120122.M
Title : SVOA CALIBRATION

Signal : TIC: BP012954.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.152	667	674	685	rVB	59354	118423	0.15%	0.042%
2	7.328	696	704	712	rBV	64156	104745	0.13%	0.038%
3	7.528	733	738	747	rVB	64430	107764	0.14%	0.039%
4	8.005	812	819	828	rBV	2163972	3413341	4.35%	1.223%
5	8.705	932	938	946	rBV	69152	119489	0.15%	0.043%
6	9.169	1009	1017	1024	rBV2	70201	121533	0.15%	0.044%
7	10.434	1225	1232	1241	rBV	83169	154311	0.20%	0.055%
8	10.822	1290	1298	1306	rBV	3270509	5511756	7.02%	1.974%
9	13.075	1674	1681	1690	rBV	70335	118858	0.15%	0.043%
10	14.051	1841	1847	1854	rBV	224804	301503	0.38%	0.108%
11	14.334	1889	1895	1902	rBV	196768	304822	0.39%	0.109%
12	14.646	1940	1948	1957	rBV	5181618	7854999	10.01%	2.813%
13	14.957	1995	2001	2006	rBV	109629	158087	0.20%	0.057%
14	15.181	2034	2039	2042	rBV3	56044	91082	0.12%	0.033%
15	15.275	2042	2055	2081	rVV2	31852614	74916338	95.44%	26.833%
16	15.634	2111	2116	2122	rVB	336849	473110	0.60%	0.169%
17	16.716	2293	2300	2313	rBV2	73540	192834	0.25%	0.069%
18	17.057	2349	2358	2361	rBV2	31229049	78498638	100.00%	28.116%
19	17.181	2376	2379	2404	rVB	1878345	3672884	4.68%	1.316%
20	17.416	2410	2419	2429	rVB	6155006	8925499	11.37%	3.197%
21	17.510	2431	2435	2441	rVB	307914	438083	0.56%	0.157%
22	17.945	2504	2509	2515	rBV2	285136	430655	0.55%	0.154%
23	18.004	2516	2519	2525	rVB6	89121	127890	0.16%	0.046%
24	18.263	2559	2563	2569	rBV2	135575	187490	0.24%	0.067%
25	19.357	2745	2749	2753	rBV	237055	304501	0.39%	0.109%
26	19.728	2808	2812	2819	rVB	284101	374318	0.48%	0.134%
27	19.886	2836	2839	2843	rVB	188551	210799	0.27%	0.076%
28	20.175	2884	2888	2894	rBV	504680	632835	0.81%	0.227%
29	20.622	2961	2964	2968	rVB	165529	178118	0.23%	0.064%
30	21.480	3105	3110	3118	rBV	4850793	6535695	8.33%	2.341%
31	22.569	3291	3295	3303	rBV	1403161	2445583	3.12%	0.876%
32	22.780	3327	3331	3338	rVB2	807510	1293291	1.65%	0.463%
33	22.975	3361	3364	3373	rVB2	504918	796590	1.01%	0.285%
34	23.939	3522	3528	3534	rBV	8670763	17347001	22.10%	6.213%
35	23.998	3534	3538	3546	rVB	3297973	6746940	8.59%	2.417%
36	24.086	3547	3553	3558	rBV	3716146	7372503	9.39%	2.641%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

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-BP120122.M
Title : SVOA CALIBRATION

37	24.139	3559	3562	3575	rVB3	1252647	3046508	3.88%	1.091%
38	24.351	3593	3598	3620	rVB3	1172374	3730035	4.75%	1.336%
39	25.568	3798	3805	3808	rBV3	2362976	5648877	7.20%	2.023%
40	25.621	3809	3814	3829	rVB2	4196276	10425018	13.28%	3.734%
41	25.792	3837	3843	3848	rBV	1482312	3934413	5.01%	1.409%
42	25.886	3850	3859	3890	rVB3	5179267	21829546	27.81%	7.819%

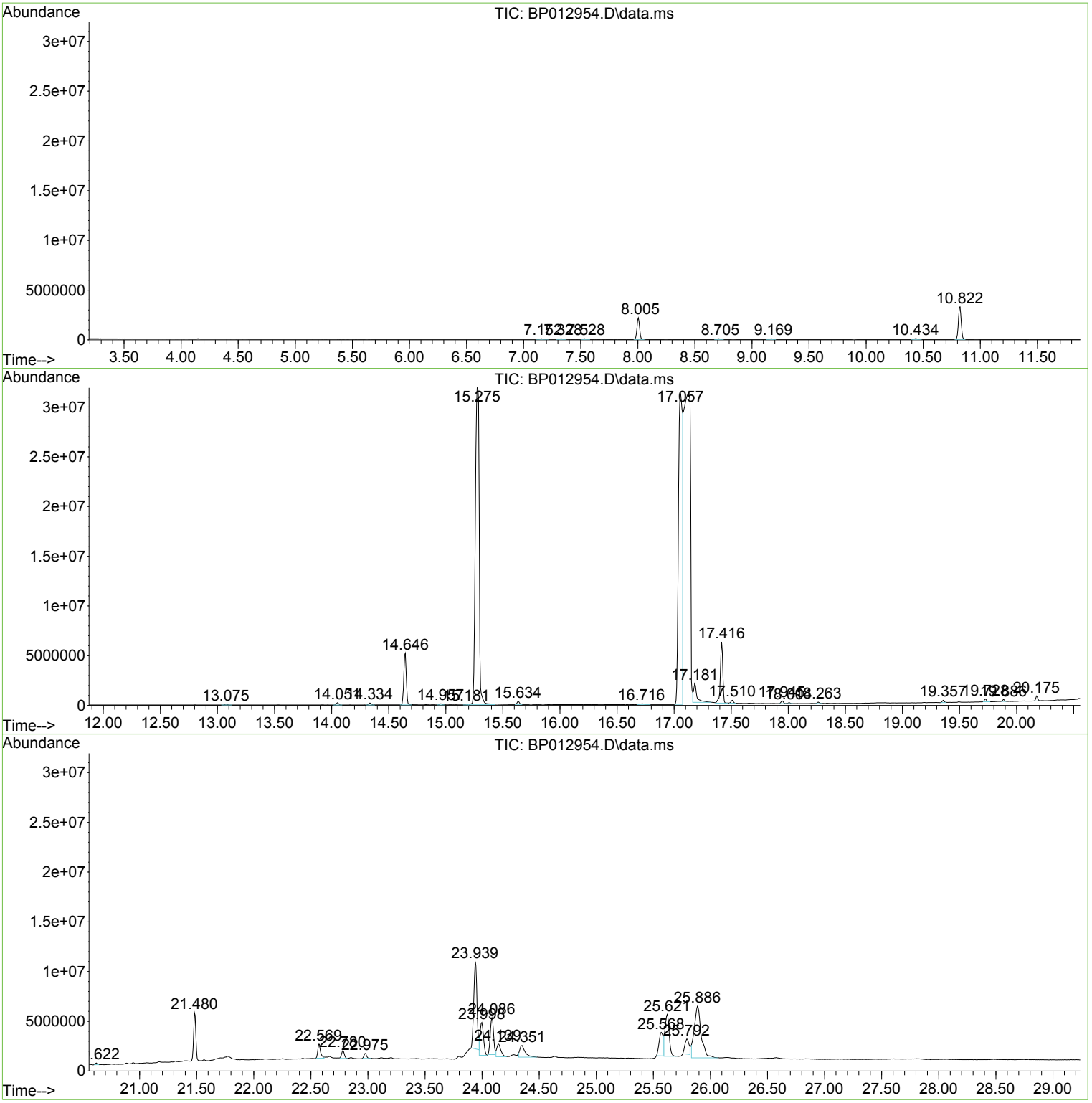
Sum of corrected areas: 279196705

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

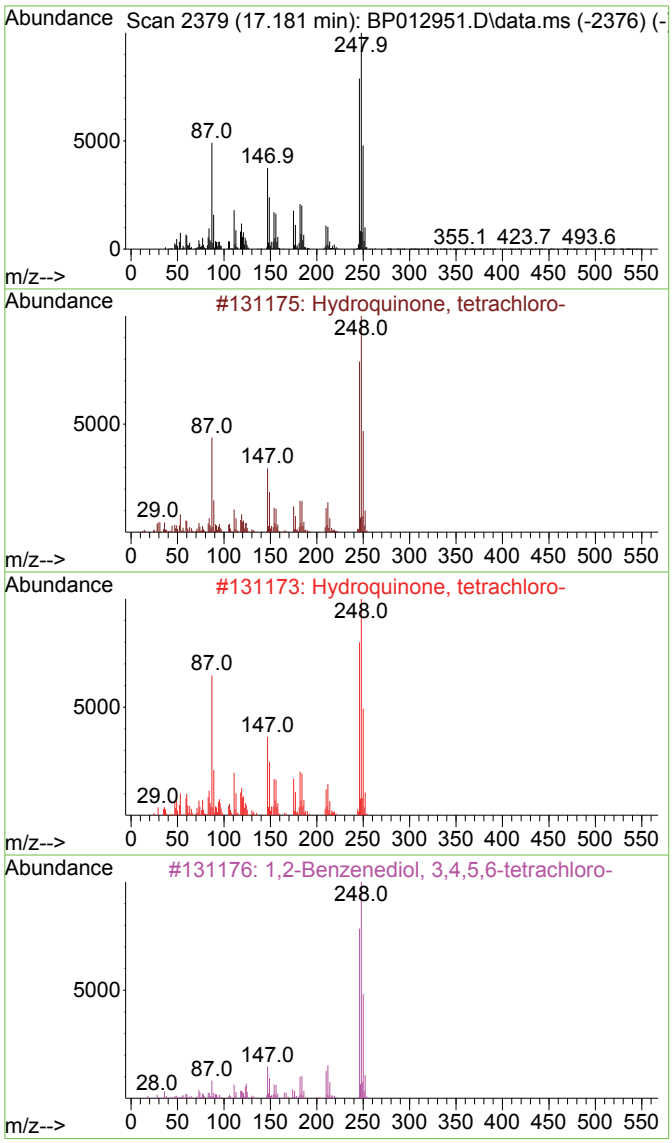
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

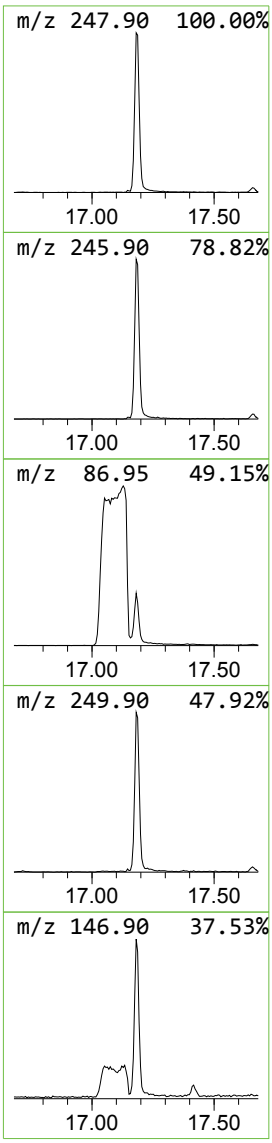
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Hydroquinone, tetrachloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.181	8.23 ng/ul	3672880	Phenanthrene-d10	17.416

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	94
2	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	94
3	1,2-Benzenediol, 3,4,5,6-tetrach...	246	C6H2Cl4O2	001198-55-6	93
4	Hydroquinone, tetrachloro-	246	C6H2Cl4O2	000087-87-6	90
5	5-Amino-4,6,8-trichloroquinoline	246	C9H5Cl3N2	1000252-19-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

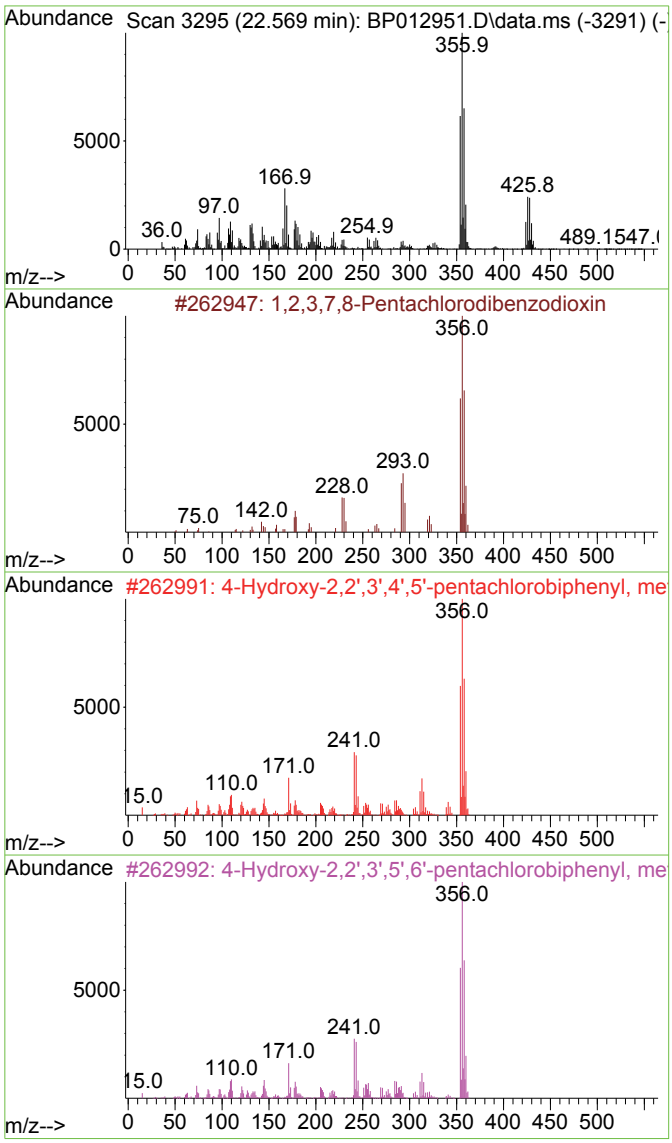
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

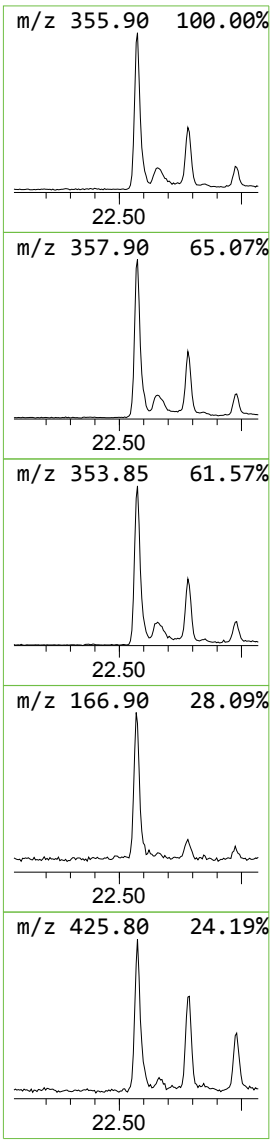
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1,2,3,7,8-Pentachlorodibenz... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.569	7.48 ng/ul	2445580	Chrysene-d12	21.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,7,8-Pentachlorodibenzodioxin	354	C12H3Cl5O2	040321-76-4	86
2	4-Hydroxy-2,2',3',4',5'-pentachl...	354	C13H7Cl5O	1000447-88-3	70
3	4-Hydroxy-2,2',3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-4	70
4	4-Hydroxy-2',3,3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-6	70
5	4-Hydroxy-2',3,4',5,6'-pentachlo...	354	C13H7Cl5O	1000447-88-7	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

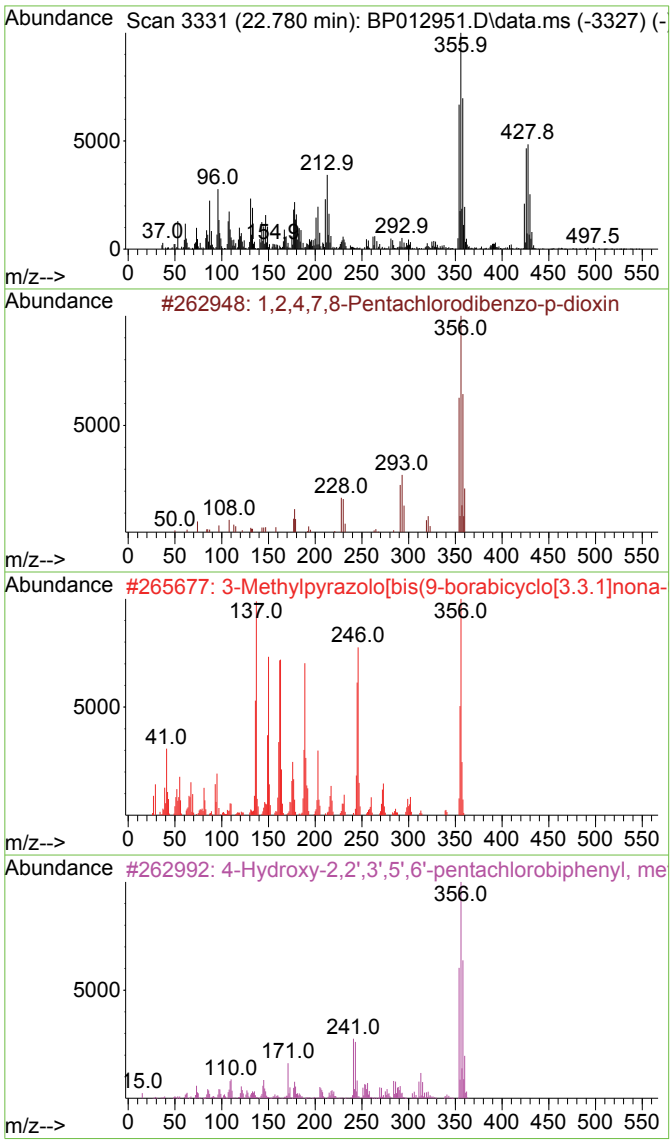
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

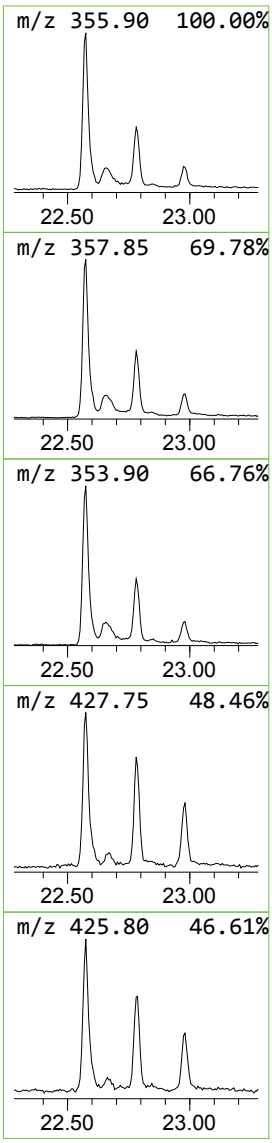
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,2,4,7,8-Pentachlorodibenz... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.780	3.83 ng/ul	1293290	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,7,8-Pentachlorodibenzo-p-d...	354	C12H3Cl5O2	058802-08-7	64
2	3-Methylpyrazolo[bis(9-borabicyc...	356	C20H34B2N2S	1000159-71-8	56
3	4-Hydroxy-2,2',3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-4	55
4	4-Hydroxy-2,2',3',4',5'-pentachl...	354	C13H7Cl5O	1000447-88-3	55
5	4-Hydroxy-2',3,3',5',6'-pentachl...	354	C13H7Cl5O	1000447-88-6	55





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

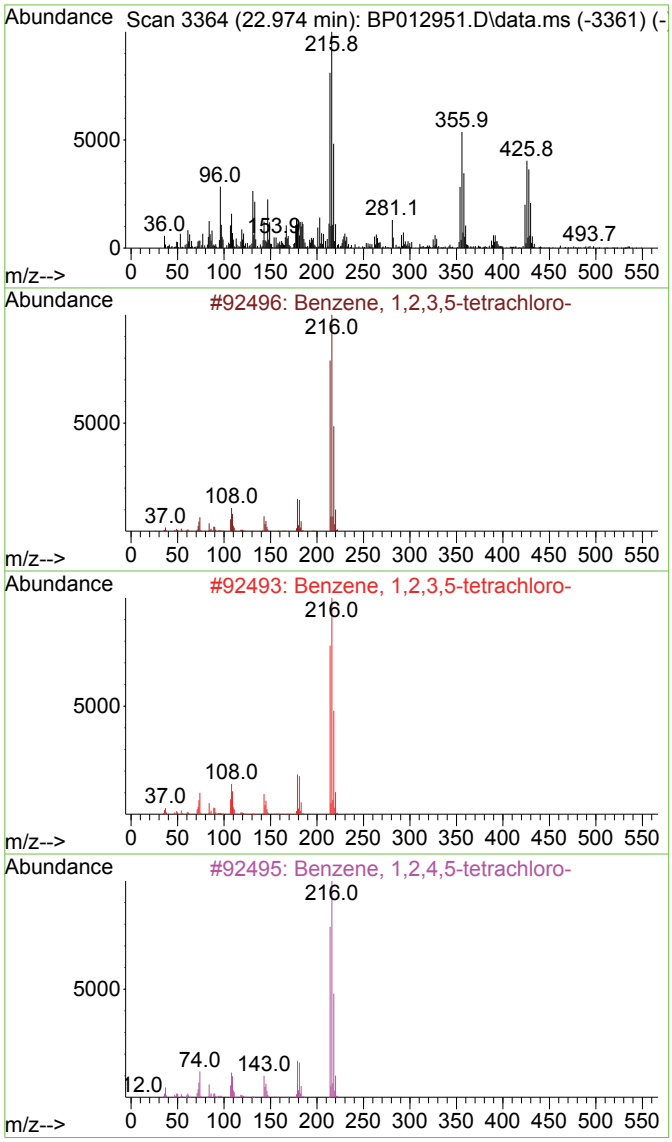
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

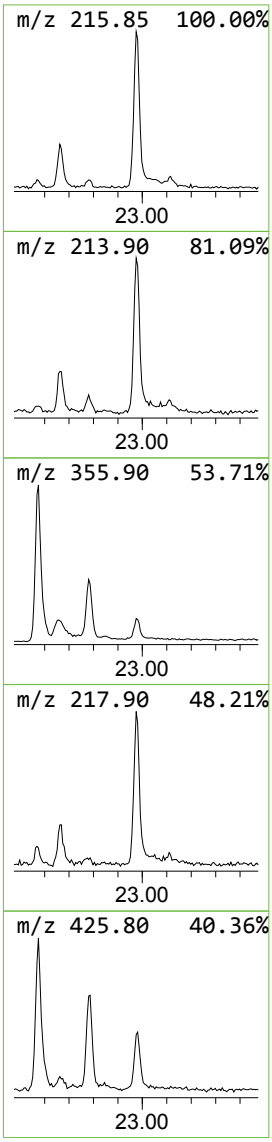
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3,5-tetrachloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.974	2.36 ng/ul	796590	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	64
2	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	64
3	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	50
4	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	50
5	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

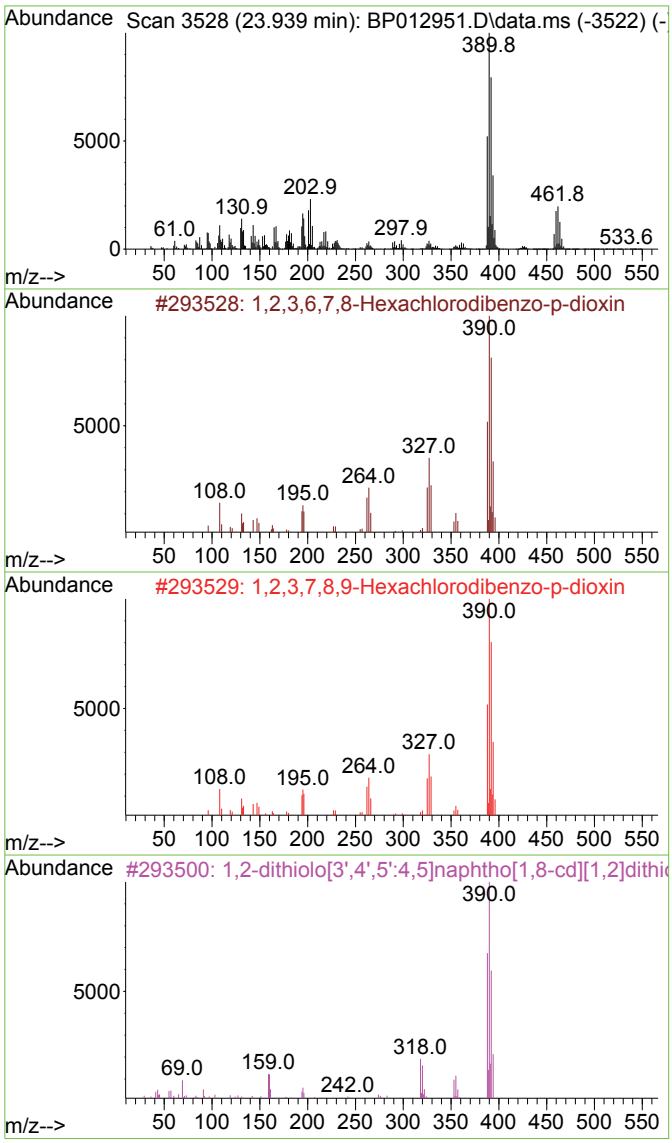
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

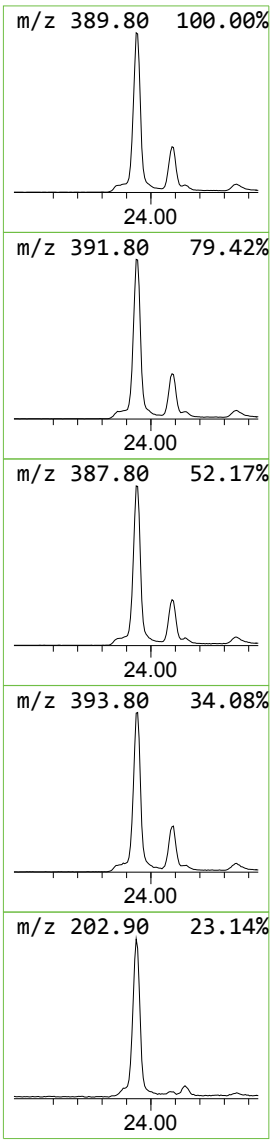
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,2,3,6,7,8-Hexachlorodiben... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.939	51.42 ng/ul	17347000	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	94
2	1,2,3,7,8,9-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	019408-74-3	90
3	1,2-dithiolo[3',4',5':4,5]naphth...	388	C10Cl4S4	1000398-83-7	40
4	4-(3-Bromopropoxy)-5-methoxy-2-n...	405	C14H20BrNO6Si	1000471-26-2	25
5	1-(4-(4-chlorophenyl)piperazin-1...	392	C14H12ClF7N2O	1010455-17-3	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
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Instrument :
BNA_P
ClientSampleId :
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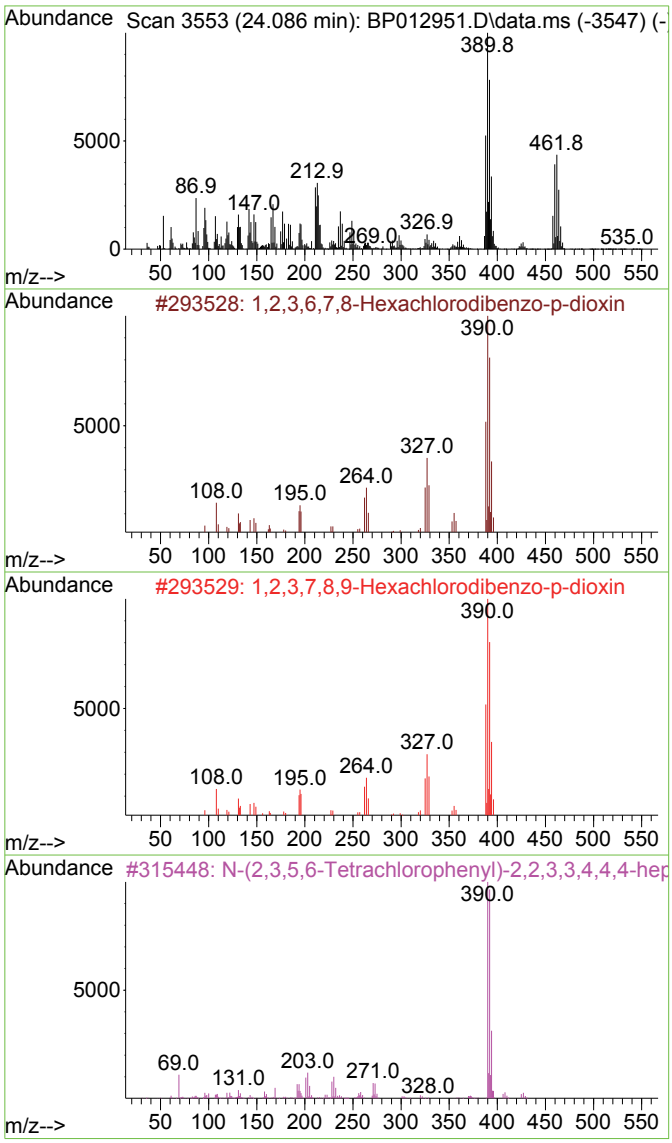
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

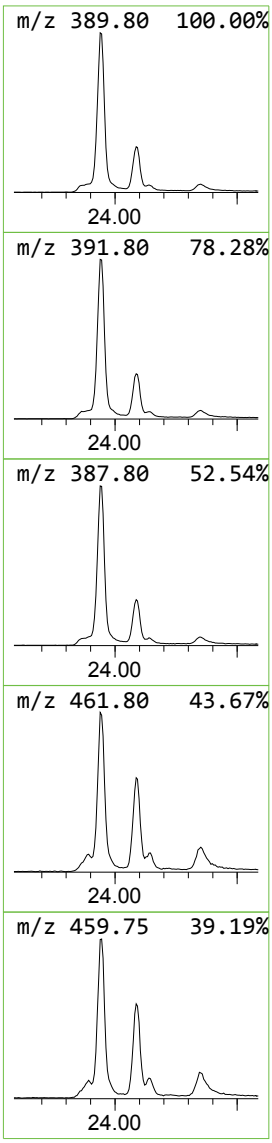
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1,2,3,7,8,9-Hexachlorodiben... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.086	21.85 ng/ul	7372500	Perylene-d12	23.998

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	53
3	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	45
4	3-Bromo-4-hydroxy-5-nitrobenzoic...	405	C13H20BrNO5Si2	1000512-83-9	38
5	1,2-dithiolo[3',4',5':4,5]naphth...	388	C10Cl4S4	1000398-83-7	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

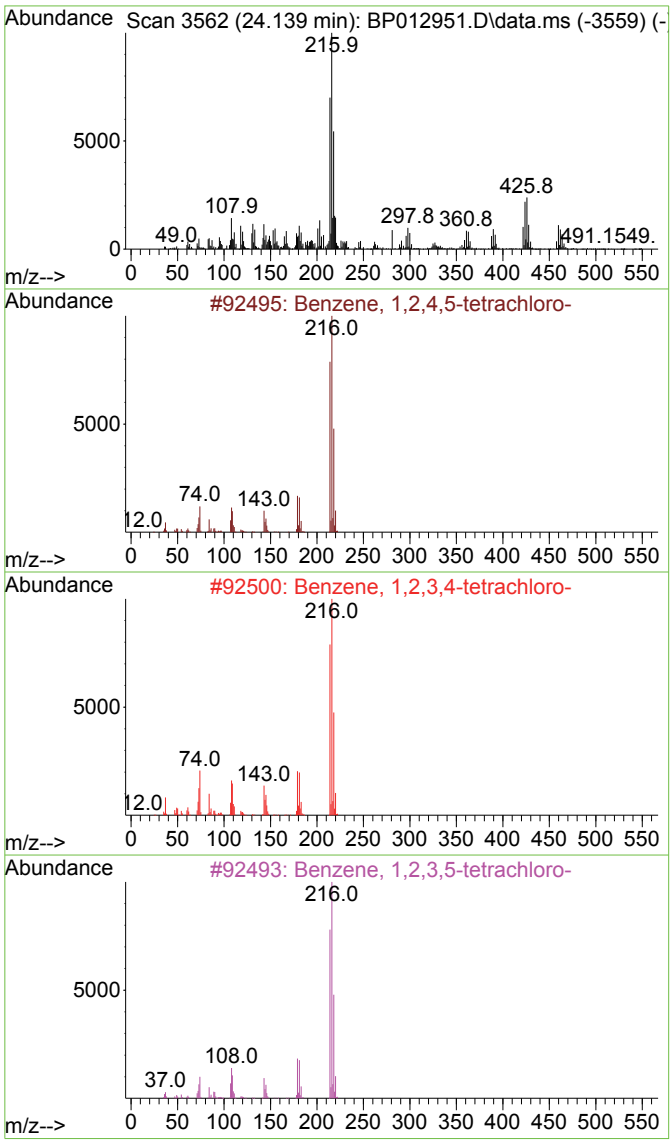
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

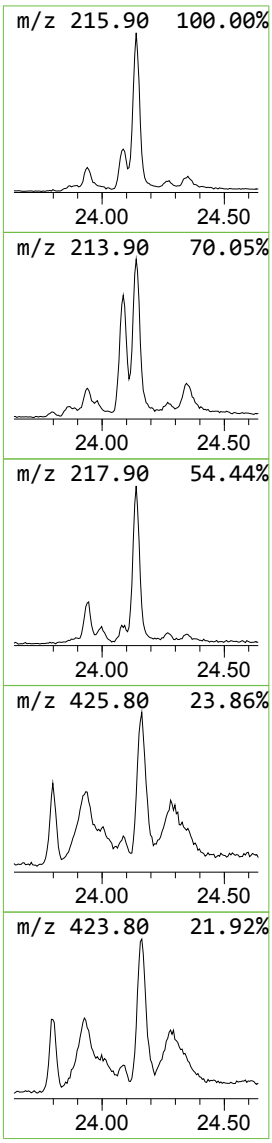
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,4,5-tetrachloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.139	9.03 ng/ul	3046510	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	70
2	Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	64
3	Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	64
4	Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	64
5	Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

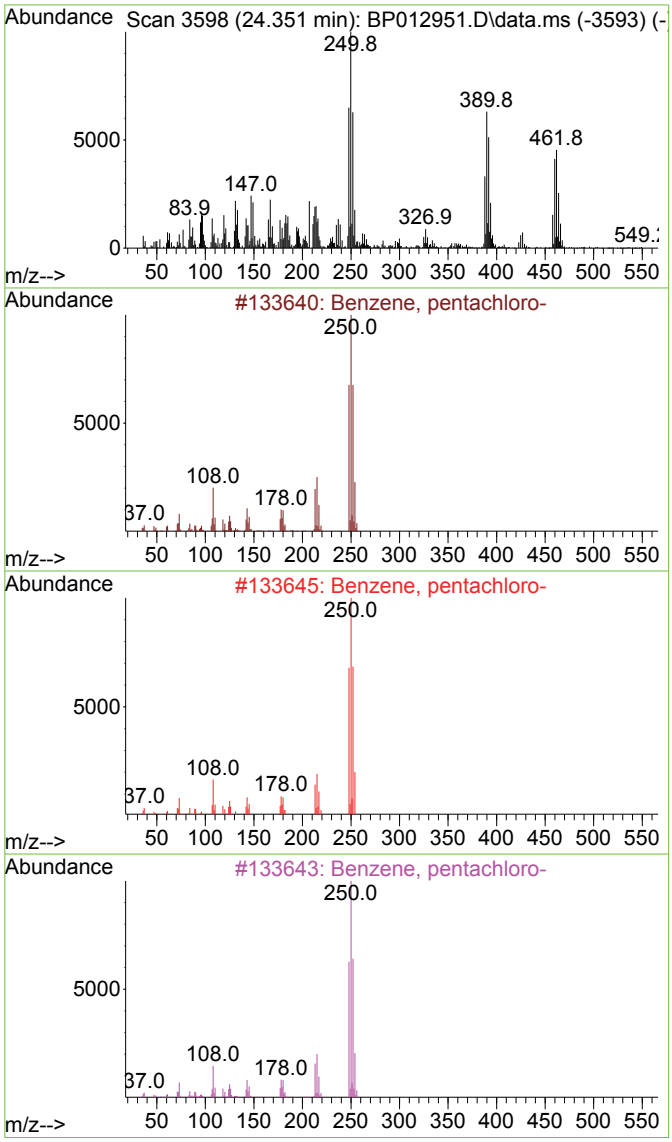
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

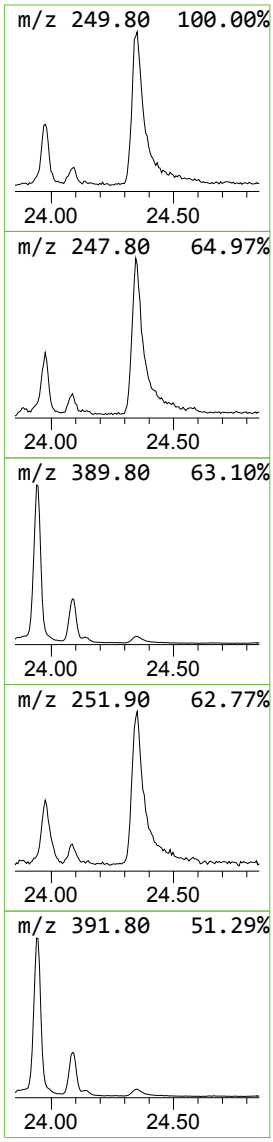
****

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.351	11.06 ng/ul	3730040	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentachloro-	248	C6HCl5	000608-93-5	46
2	Benzene, pentachloro-	248	C6HCl5	000608-93-5	41
3	Benzene, pentachloro-	248	C6HCl5	000608-93-5	38
4	Benzene, pentachloro-	248	C6HCl5	000608-93-5	35
5	Benzene, pentachloro-	248	C6HCl5	000608-93-5	35





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Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

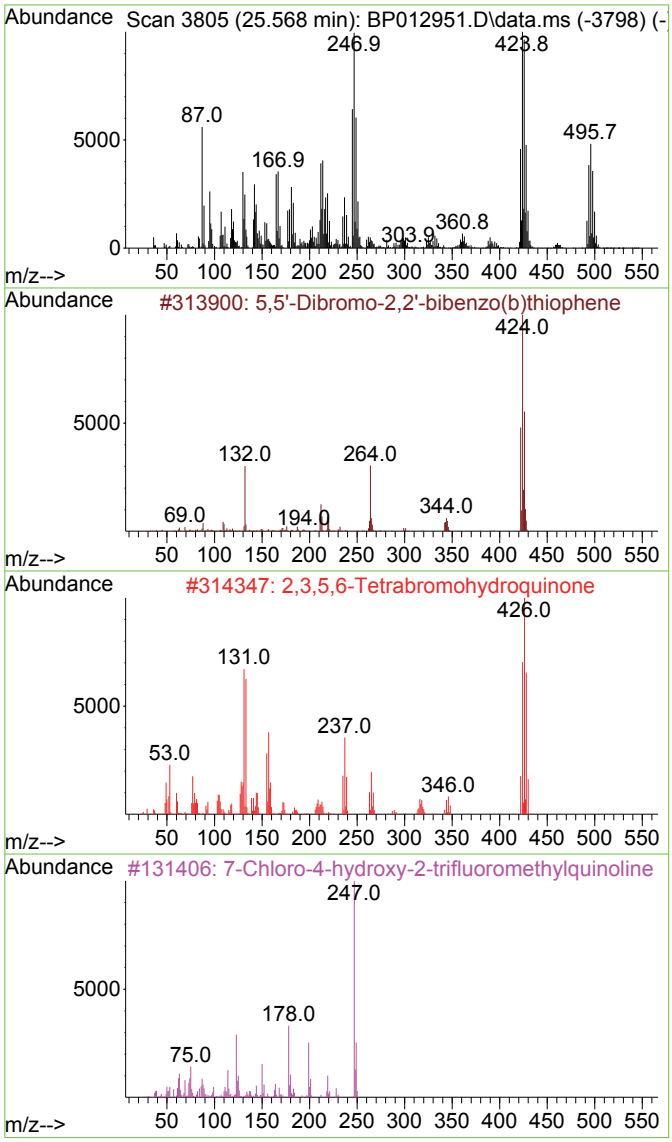
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

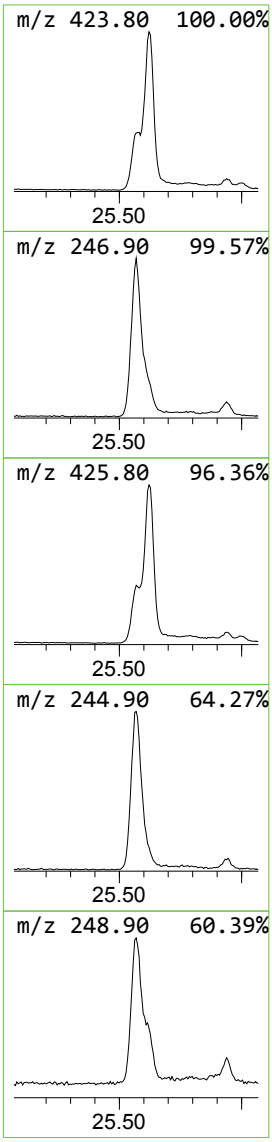
****

Peak Number 9 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.569	16.75 ng/ul	5648880	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	20
2	2,3,5,6-Tetrabromohydroquinone	422	C6H2Br4O2	002641-89-6	14
3	7-Chloro-4-hydroxy-2-trifluorome...	247	C10H5ClF3NO	057124-20-6	11
4	Acetic acid, (4-chloro-2-methylp...	424	C25H41ClO3	1000415-16-9	10
5	Tetrabromocatechol	422	C6H2Br4O2	000488-47-1	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

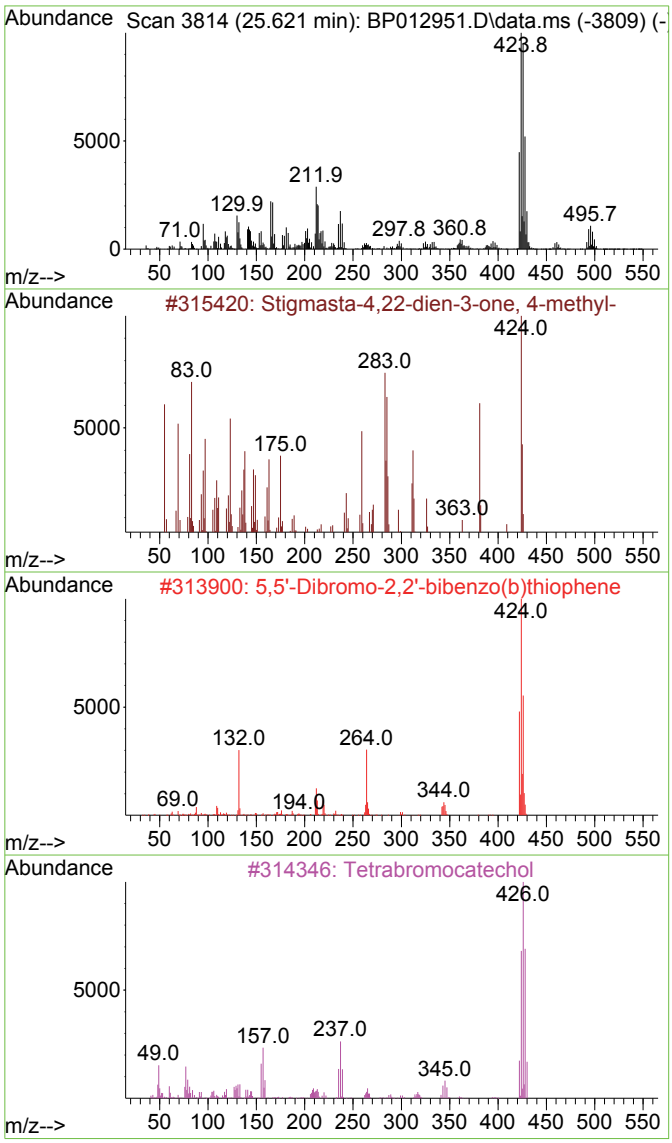
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

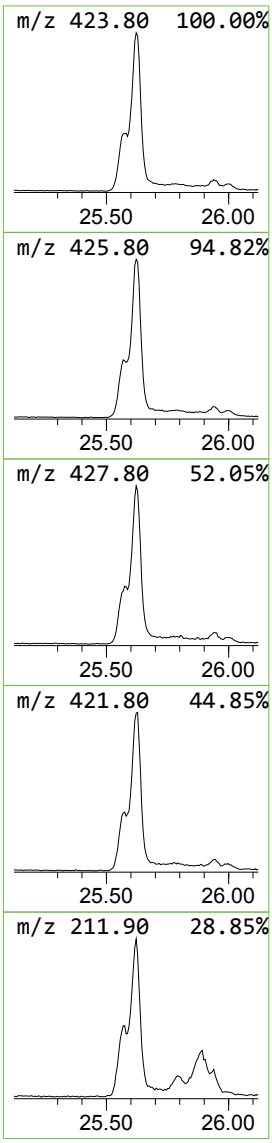
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Stigmasta-4,22-dien-3-one, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.621	30.90 ng/ul	10425000	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmasta-4,22-dien-3-one, 4-met...	424	C30H48O	057884-74-9	68
2	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	53
3	Tetrabromocatechol	422	C6H2Br4O2	000488-47-1	49
4	Phenol, 2,3,4,5-tetrabromo-6-met...	420	C7H4Br4O	000576-55-6	47
5	3,4,5-Tribromophenol, trifluoroa...	424	C8H2Br3F3O2	1000448-03-8	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

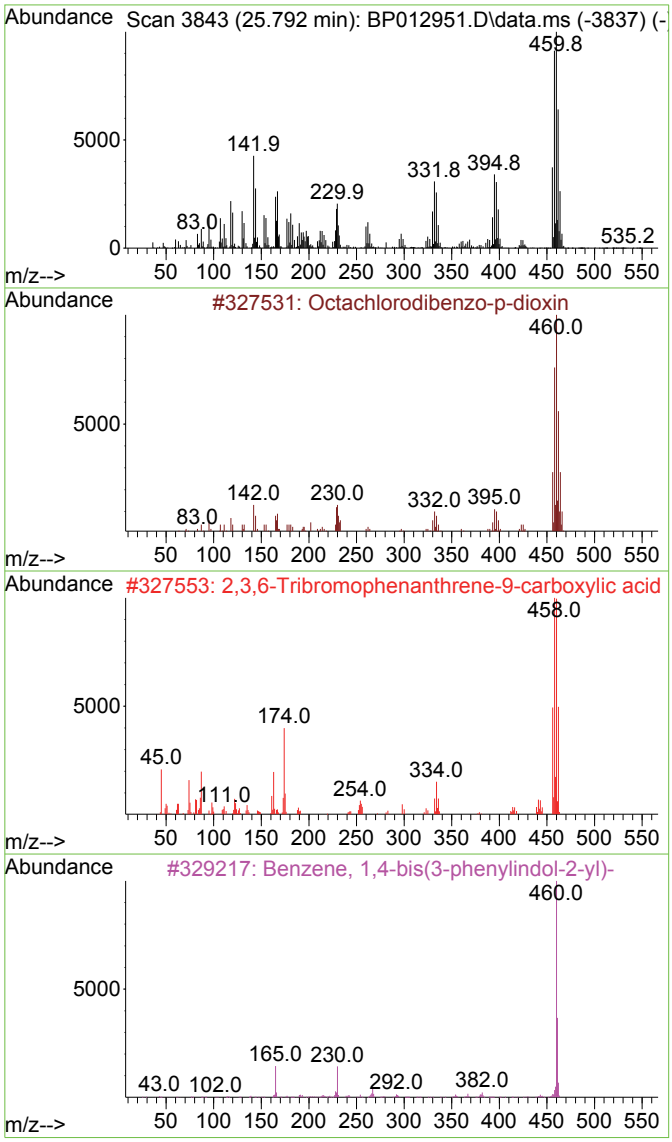
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

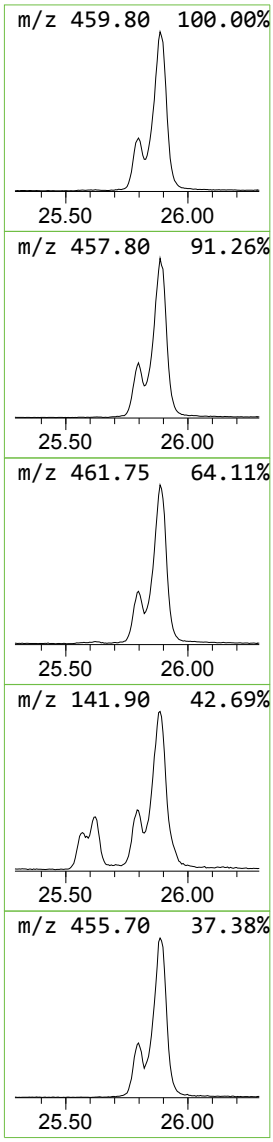
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Octachlorodibenzo-p-dioxin Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.792	11.66 ng/ul	3934410	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	95
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	43
3	Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	14
4	3-Chloro-.alpha.-[2-pyridyl]-2,4...	458	C29H31Cl3N2O	035158-92-0	14
5	1,1':4',1'':4'',1''':4''',1'''':...	458	C36H26	004499-83-6	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

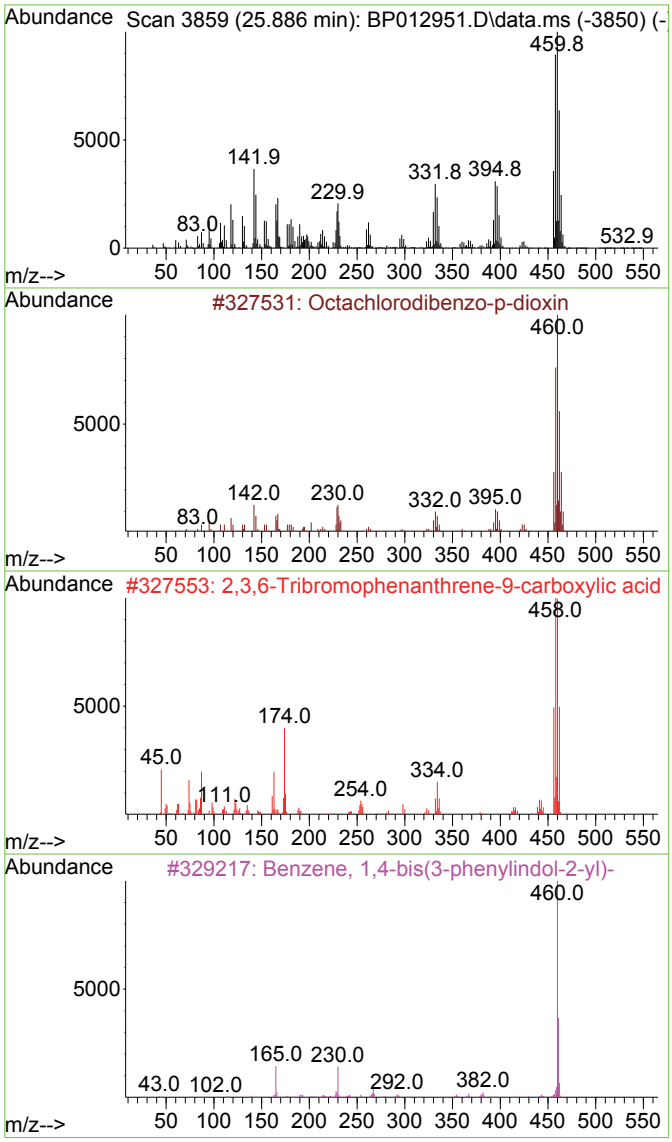
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
Quant Title : SVOA CALIBRATION

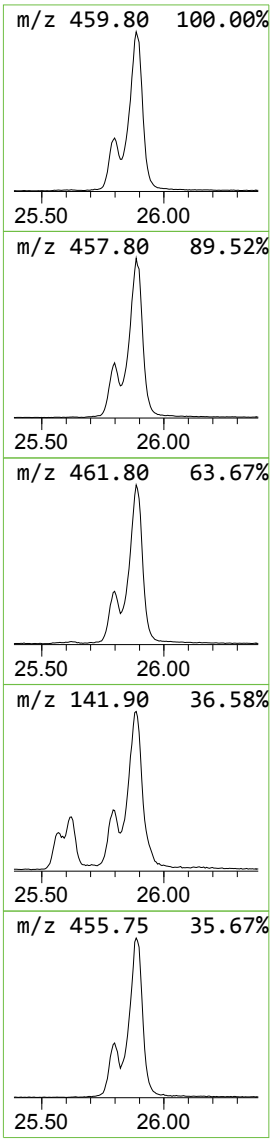
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.886	64.71 ng/ul	21829500	Perylene-d12	23.998

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	87
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	43
3	Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10
4	Cholest-8(14)-en-3-ol, (3.beta.,...	458	C30H54OSi	055282-50-3	10
5	Cholest-7-en-3-ol, (3.beta.,5.al...	458	C30H54OSi	002665-03-4	10





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
Data File : BP012951.D
Acq On : 05 Dec 2022 14:04
Operator : CG/JU
Sample : N5738-15DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBNJ4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.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
Hydroquinone, t...	17.181	8.2	ng/ul	3672880	4	17.416	8925500	20.0
1,2,3,7,8-Penta...	22.569	7.5	ng/ul	2445580	5	21.480	6535700	20.0
1,2,4,7,8-Penta...	22.780	3.8	ng/ul	1293290	6	23.998	6746940	20.0
Benzene, 1,2,3,...	22.974	2.4	ng/ul	796590	6	23.998	6746940	20.0
1,2,3,6,7,8-Hex...	23.939	51.4	ng/ul	17347000	6	23.998	6746940	20.0
1,2,3,7,8,9-Hex...	24.086	21.9	ng/ul	7372500	6	23.998	6746940	20.0
Benzene, 1,2,4,...	24.139	9.0	ng/ul	3046510	6	23.998	6746940	20.0
unknown-01	24.351	11.1	ng/ul	3730040	6	23.998	6746940	20.0
unknown-02	25.569	16.8	ng/ul	5648880	6	23.998	6746940	20.0
Stigmasta-4,22-...	25.621	30.9	ng/ul	10425000	6	23.998	6746940	20.0
Octachlorodiben...	25.792	11.7	ng/ul	3934410	6	23.998	6746940	20.0
unknown-03	25.886	64.7	ng/ul	21829500	6	23.998	6746940	20.0