LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4477

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-BP010523.M
Title : SVOA CALIBRATION

Signal : TIC: BP013426.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	17	21	27	rVB	118314	155391	0.16%	0.058%
2	3.734	90	93	119	rBV	1427125	2223346	2.32%	0.829%
3	4.522	221	227	242	rBV3	87831	266946	0.28%	0.100%
4	5.617	406	413	416	rBV3	51097	124959	0.13%	0.047%
5	6.452	546	555	565	rBV2	38682	132460	0.14%	0.049%
6	7.075	648	661	674	rBV	1825397	2783082	2.90%	1.038%
7	7.246	683	690	702	rVB	1437721	2230247	2.32%	0.832%
8	7.446	717	724	733	rBV	1746218	2770207	2.89%	1.033%
9	7.916	796	804	810	rBV	1714300	2656036	2.77%	0.991%
10	8.328	868	874	882	rBV	42201	99137	0.10%	0.037%
11	8.622	917	924	935	rBV	1982130	3213915	3.35%	1.199%
12	9.081	995	1002	1017	rBV	1573997	2549228	2.66%	0.951%
13	9.805	1118	1125	1134	rBV	1280893	2100319	2.19%	0.783%
14	10.346	1209	1217	1229	rBV	2070944	3366243	3.51%	1.256%
15	10.728	1274	1282	1290	rVB	2303724	3795442	3.96%	1.416%
16	10.869	1298	1306	1321	rBV	1775829	3200511	3.34%	1.194%
17	12.910	1648	1653	1657	rBV	122869	161328	0.17%	0.060%
18	12.969	1657	1663	1668	rVB	162684	228409	0.24%	0.085%
19	13.204	1699	1703	1707	rBV	96092	126110	0.13%	0.047%
20	13.969	1825	1833	1842	rBV	3473535	4598588	4.79%	1.715%
21	14.251	1875	1881	1896	rBV	3665453	5473903	5.71%	2.042%
22	14.563	1923	1934	1944	rBV	3486533	5203719	5.42%	1.941%
23	14.763	1961	1968	1986	rBV	1275553	1989710	2.07%	0.742%
24	15.510	2089	2095	2097	rBV	100729	151471	0.16%	0.056%
25	15.557	2097	2103	2108	rVV	5087407	7135526	7.44%	2.662%
26	15.604	2108	2111	2115	rVV	72206	97072	0.10%	0.036%
27	15.681	2115	2124	2132	rVV	1405076	1922947	2.00%	0.717%
28	15.840	2141	2151	2158	rVV	20854946	29126944	30.36%	10.864%
29	15.922	2158	2165	2169	rVB2	128599	243043	0.25%	0.091%
30	16.151	2198	2204	2210	rBV	15129846	19824792	20.66%	7.395%
31	16.204	2210	2213	2218	rVB	111617	142202	0.15%	0.053%
32	16.375	2237	2242	2247	rBV	478817	611749	0.64%	0.228%
33	16.457	2252	2256	2261	rVB2	156690	221994	0.23%	0.083%
34	16.545	2265	2271	2278	rVB2	255110	401986	0.42%	0.150%
35	16.681	2290	2294	2302	rVB3	75081	107899	0.11%	0.040%
36	16.845	2318	2322	2327	rVB	90623	120501	0.13%	0.045%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

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-BP010523.M
Title : SVOA CALIBRATION

37	17.192	2376	2381	2385	rBV	701271	948078	0.99%	0.354%
38	17.228	2385	2387	2392	rVB	274263	330347	0.34%	0.123%
39	17.322	2395	2403	2413	rBV	4371340	6402283	6.67%	2.388%
40	17.416	2413	2419	2427	rVV	4958028	7367493	7.68%	2.748%
41	17.492	2427	2432	2444	rVB	478494	780719	0.81%	0.291%
42	17.769	2475	2479	2482	rBV	126083	157438	0.16%	0.059%
43	17.910	2497	2503	2510	rBV	889904	1740259	1.81%	0.649%
44	18.039	2518	2525	2530	rBV3	383458	647341	0.67%	0.241%
45	18.157	2541	2545	2547	rBV	180859	226043	0.24%	0.084%
46	18.192	2547	2551	2559	rVB2	651924	960910	1.00%	0.358%
47	18.369	2576	2581	2587	rVV	670871	862590	0.90%	0.322%
48	18.428	2587	2591	2594	rVV	528564	625091	0.65%	0.233%
49	18.469	2594	2598	2602	rVB	399721	460619	0.48%	0.172%
50	18.645	2623	2628	2633	rVV	671497	911742	0.95%	0.340%
51	18.686	2633	2635	2640	rVV	211254	260406	0.27%	0.097%
52	18.745	2641	2645	2648	rVV	173252	241327	0.25%	0.090%
53	18.781	2648	2651	2653	rVV	188175	222644	0.23%	0.083%
54	18.816	2653	2657	2662	rVB	372369	467995	0.49%	0.175%
55	18.910	2670	2673	2679	rVB4	145267	181645	0.19%	0.068%
56	19.110	2704	2707	2711	rBV	102823	125000	0.13%	0.047%
57	19.163	2712	2716	2718	rBV	216829	271122	0.28%	0.101%
58	19.198	2719	2722	2725	rVB2	169636	206307	0.22%	0.077%
59	19.422	2757	2760	2764	rVB2	191310	207455	0.22%	0.077%
60	19.651	2794	2799	2806	rBV	7172742	9169404	9.56%	3.420%
61	21.098	3041	3045	3059	rBV2	2599821	4053892	4.23%	1.512%
62	21.304	3075	3080	3085	rBV2	41875177	95947898	100.00%	35.789%
63	21.410	3094	3098	3109	rVB	5133030	7206295	7.51%	2.688%
64	23.716	3484	3490	3505	rVB2	4701450	10313295	10.75%	3.847%
65	23.869	3511	3516	3526	rVB2	3508684	7243152	7.55%	2.702%

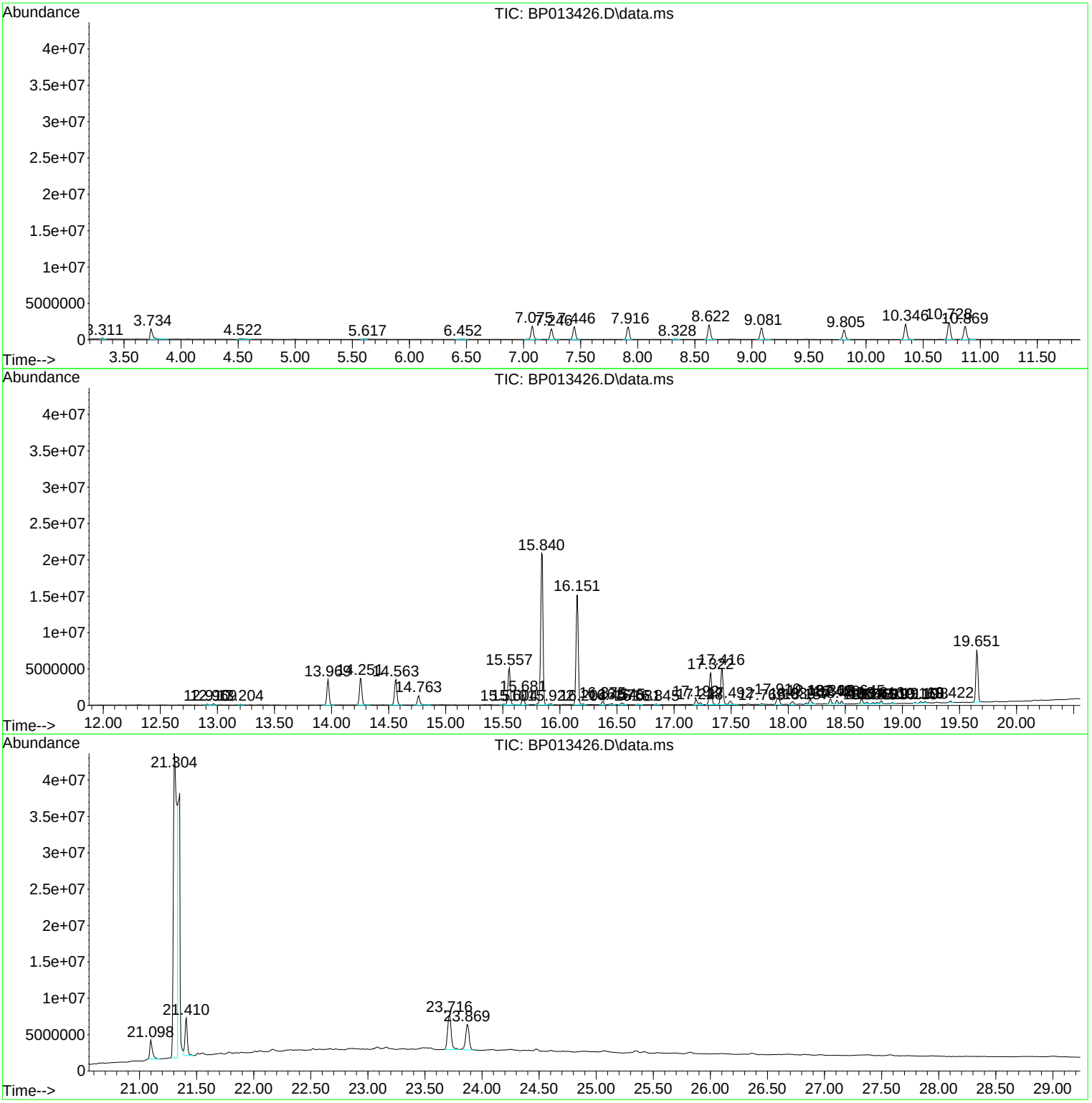
Sum of corrected areas: 268096152

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

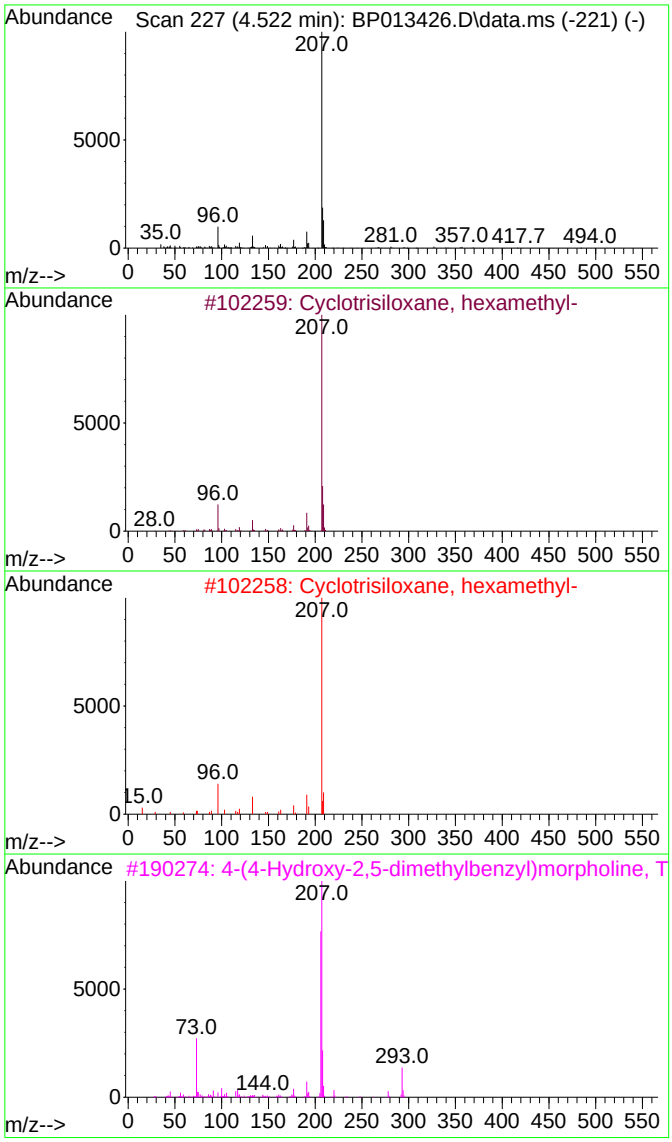
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

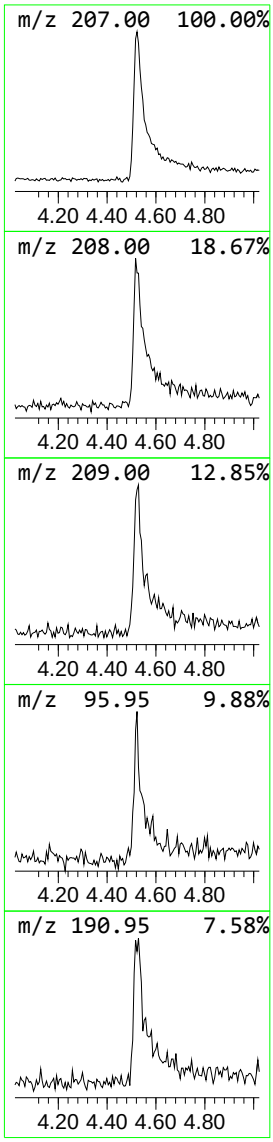
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclotrisiloxane, hexamethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.522	2.01 ng/ul	266946	1,4-Dichlorobenzene-d4	7.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	91
2	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	78
3	4-(4-Hydroxy-2,5-dimethylbenzyl)...	293	C16H27NO2Si	1000507-09-7	64
4	Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	58
5	Tetrasiloxane, decamethyl-	310	C10H30O3Si4	000141-62-8	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

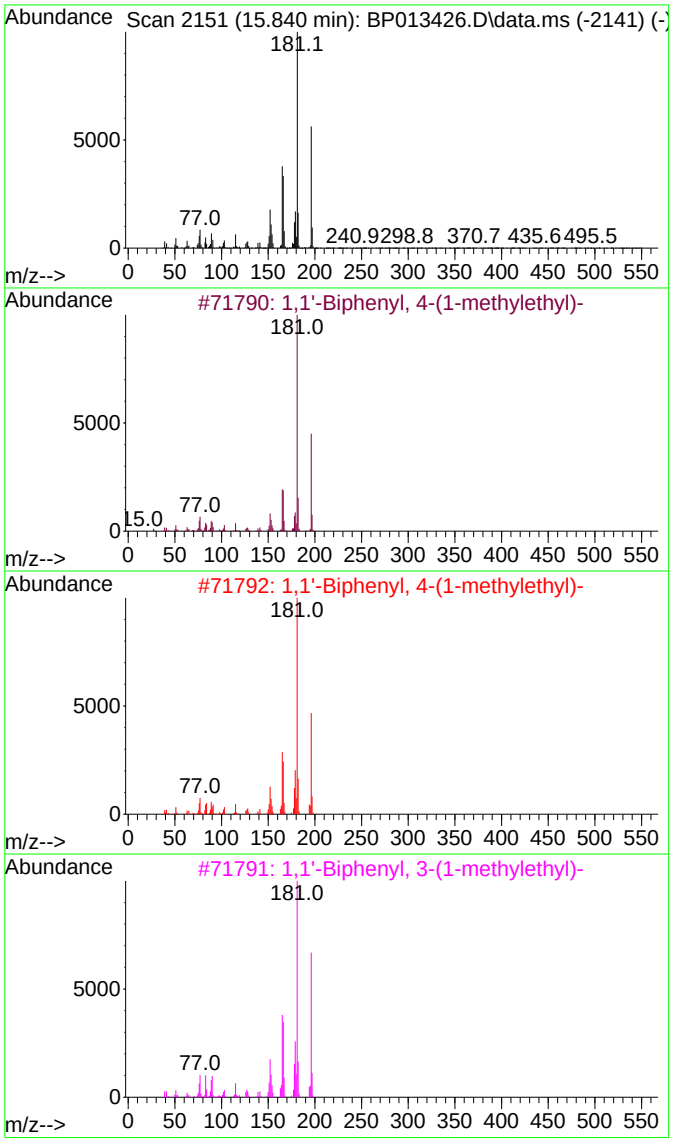
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

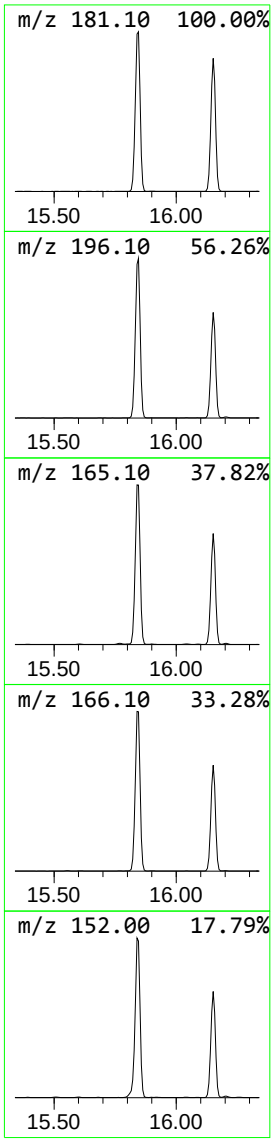
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1,1'-Biphenyl, 4-(1-methyle... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.839	111.95 ng/ul	29126900	Acenaphthene-d10	14.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	96
2	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	94
3	1,1'-Biphenyl, 3-(1-methylethyl)-	196	C15H16	020282-30-8	94
4	Benzene, 3,5-dimethyl-1-(phenylm...	196	C15H16	028122-27-2	94
5	Benzene, 2,4-dimethyl-1-(phenylm...	196	C15H16	028122-28-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

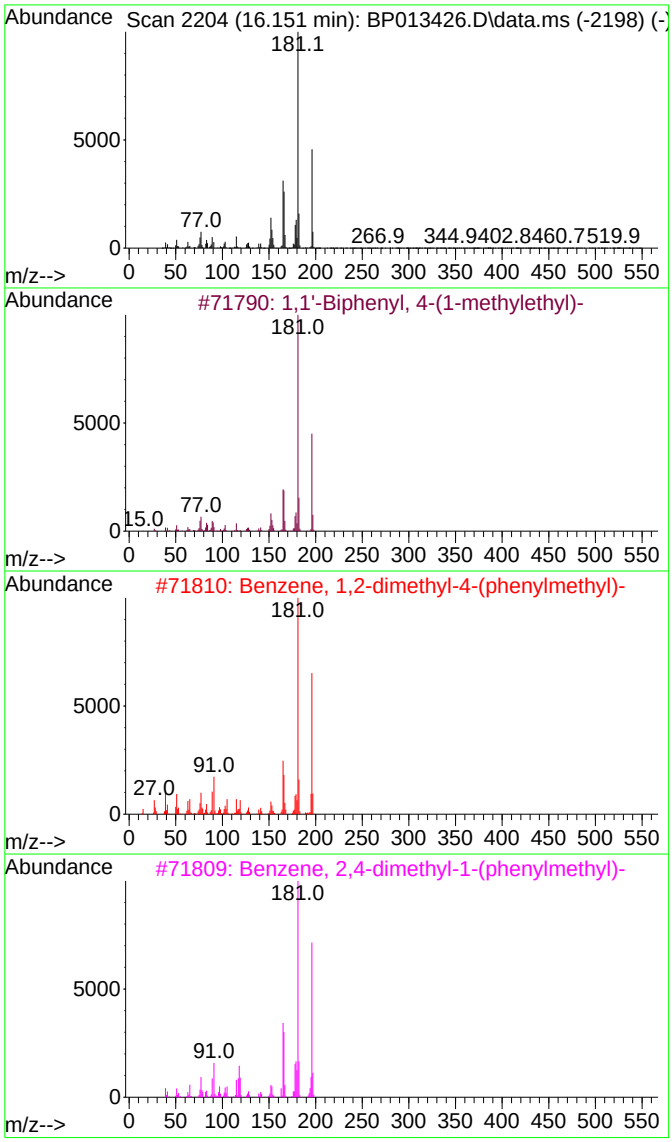
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

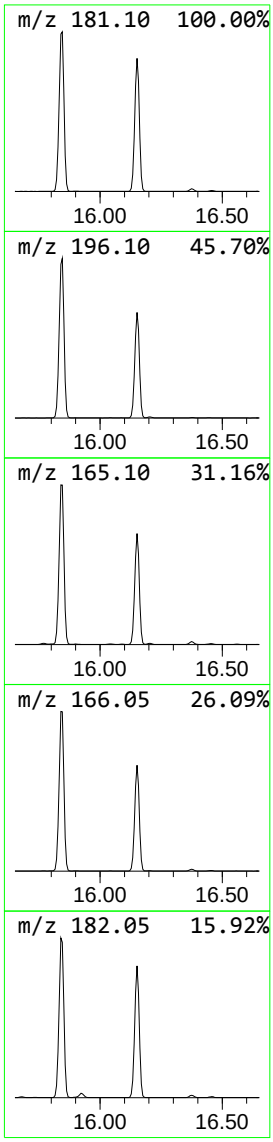
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,2-dimethyl-4-(ph... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.151	61.93 ng/ul	19824800	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	97
2	Benzene, 1,2-dimethyl-4-(phenylm...	196	C15H16	013540-56-2	94
3	Benzene, 2,4-dimethyl-1-(phenylm...	196	C15H16	028122-28-3	91
4	Methane, di-p-tolyl-	196	C15H16	004957-14-6	91
5	Methane, di-p-tolyl-	196	C15H16	004957-14-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

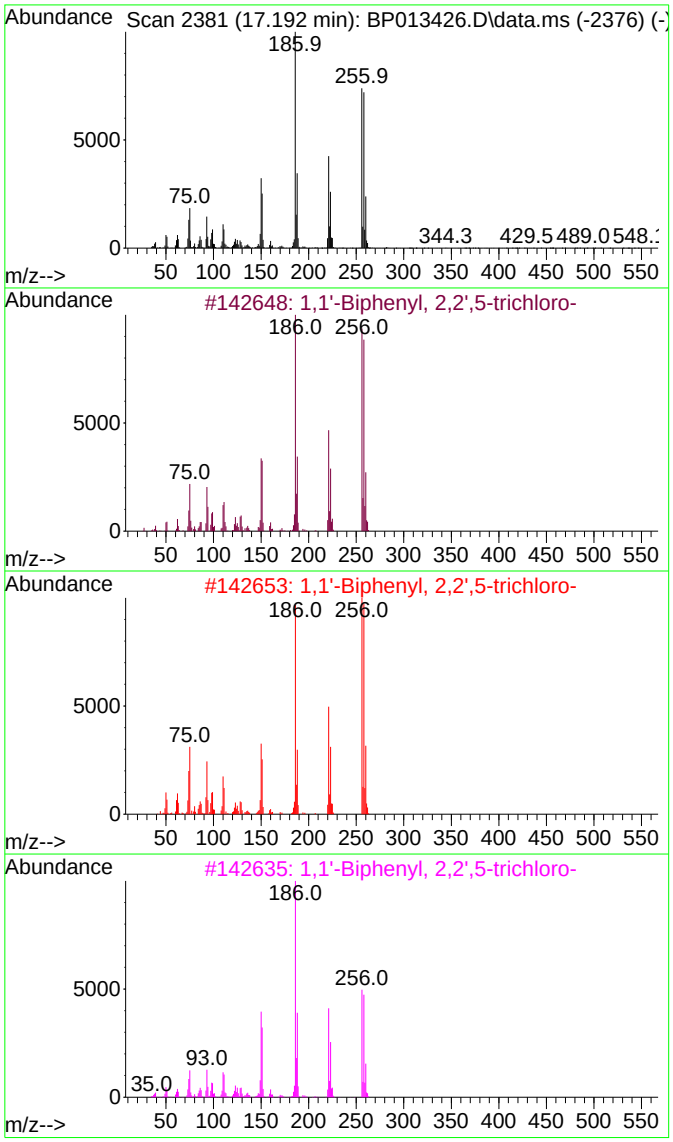
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

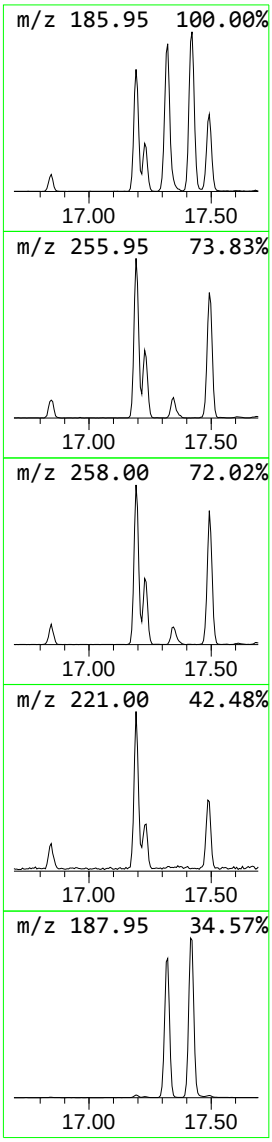
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.192	2.96 ng/ul	948078	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
4	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
5	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

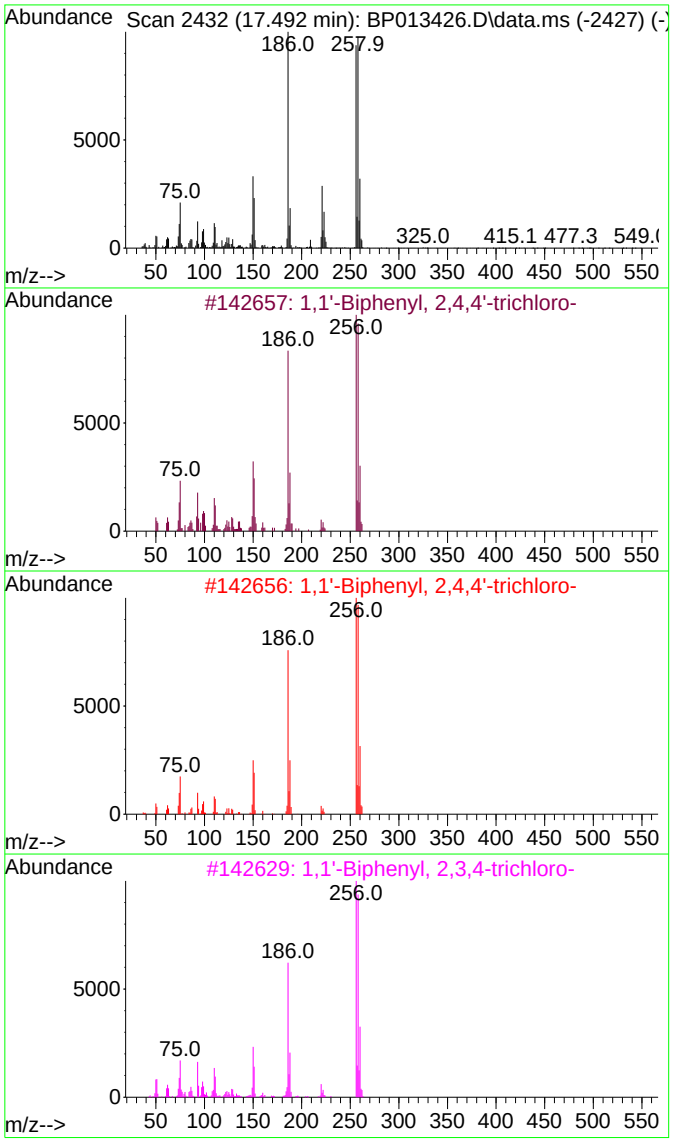
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

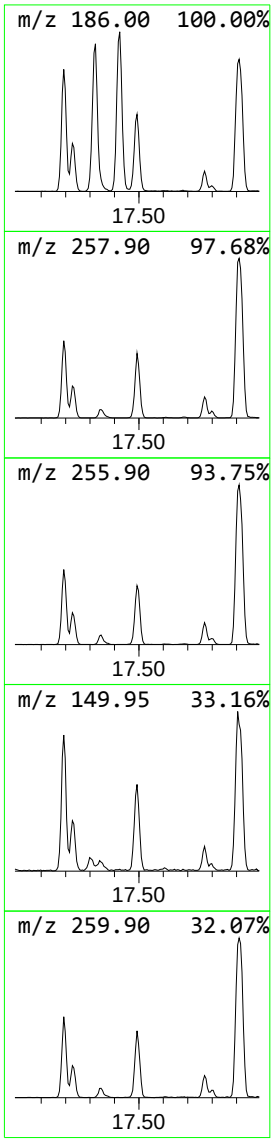
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.492	2.44 ng/ul	780719	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	95
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	94
4	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	94
5	1,1'-Biphenyl, 2,3',4-trichloro-	256	C12H7Cl3	055712-37-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

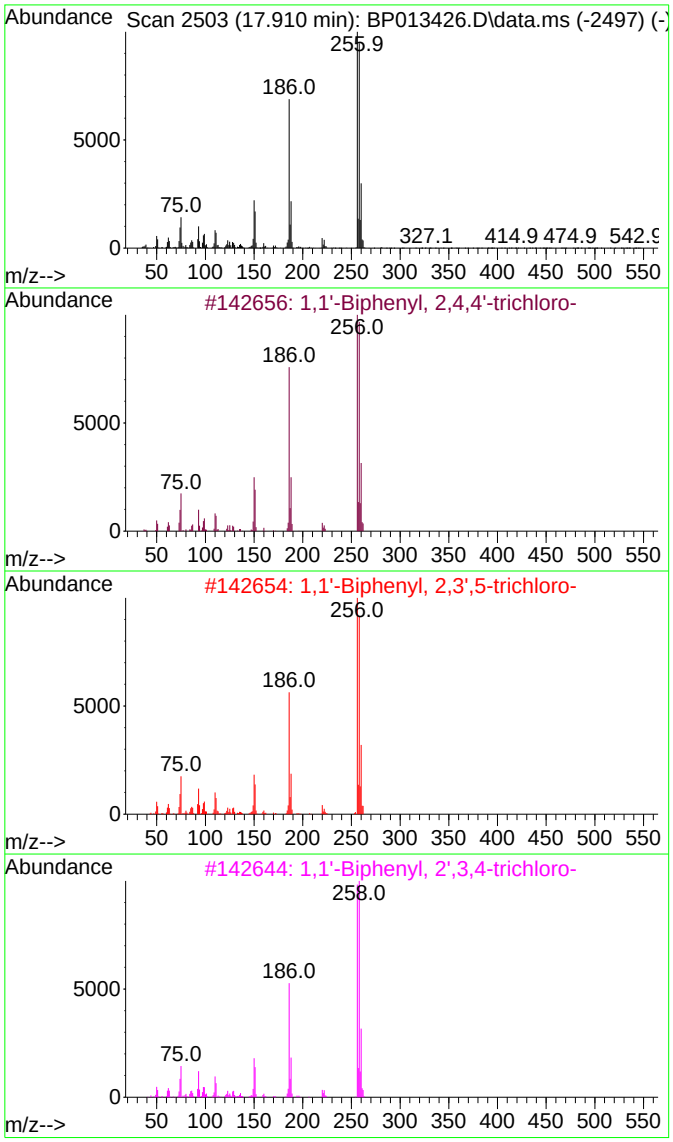
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

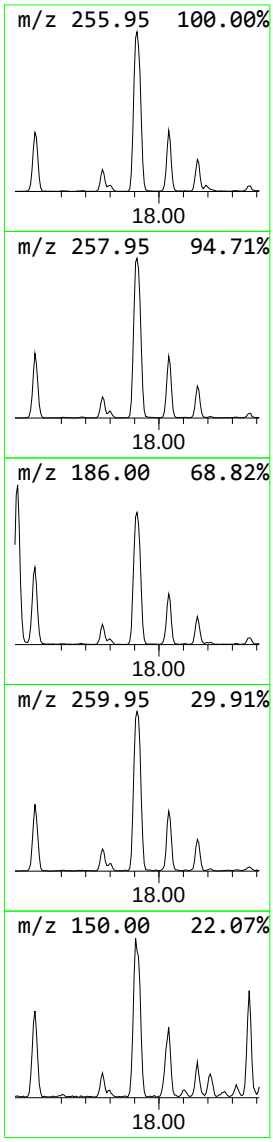
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1,1'-Biphenyl, 2,3',5-trich... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.910	5.44 ng/ul	1740260	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
3	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
4	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
5	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

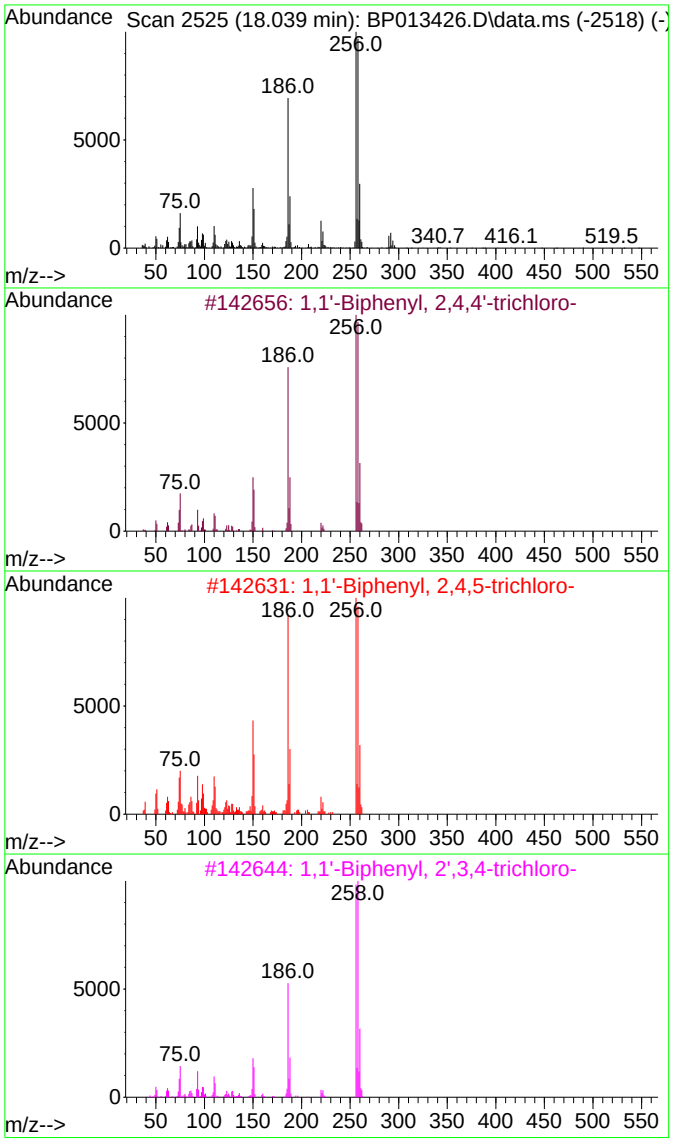
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

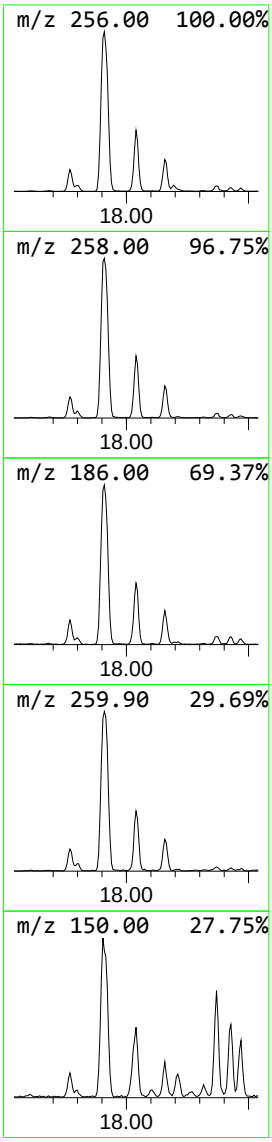
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,1'-Biphenyl, 2,4,5-trichl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	2.02 ng/ul	647341	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	99
3	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
4	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98
5	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	98





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Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

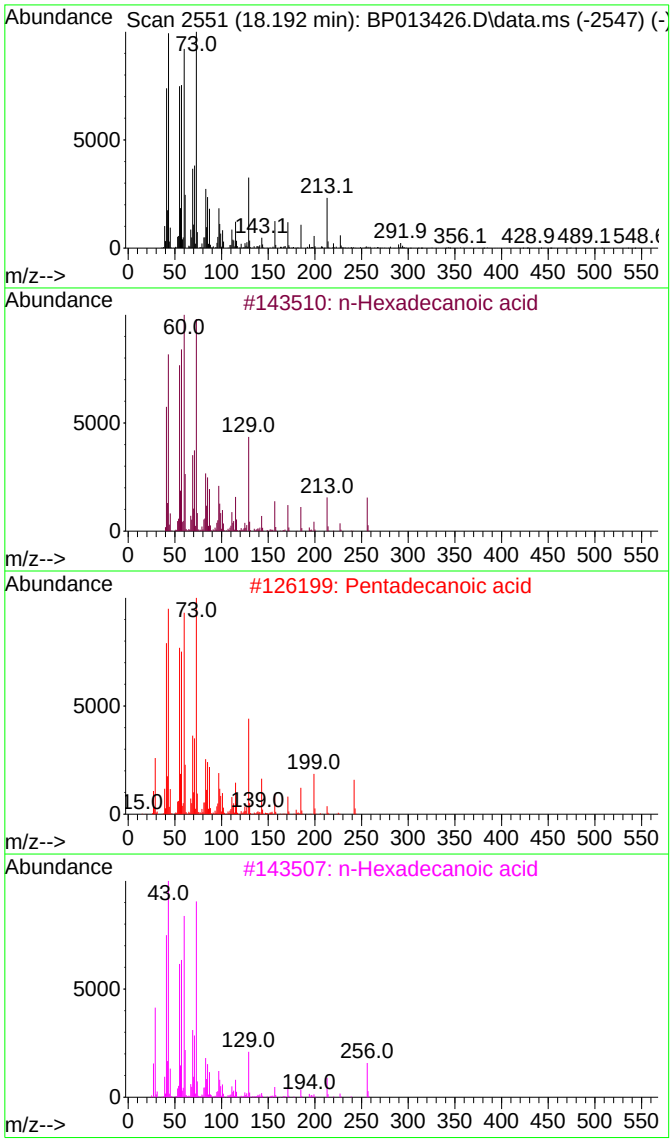
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

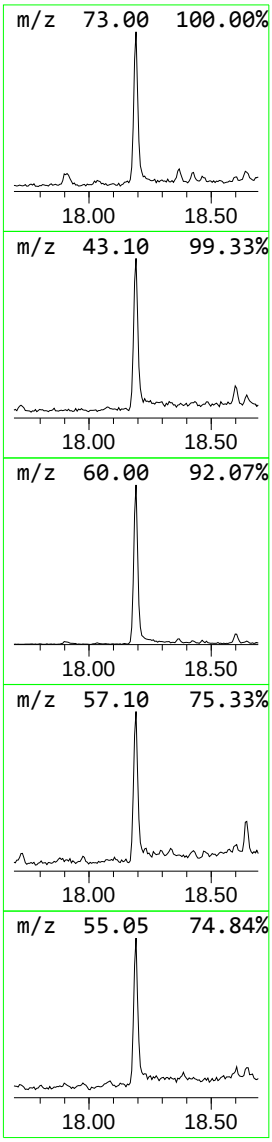
****

Peak Number 8 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	3.00 ng/ul	960910	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
4	Octadecanoic acid	284	C18H36O2	000057-11-4	81
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

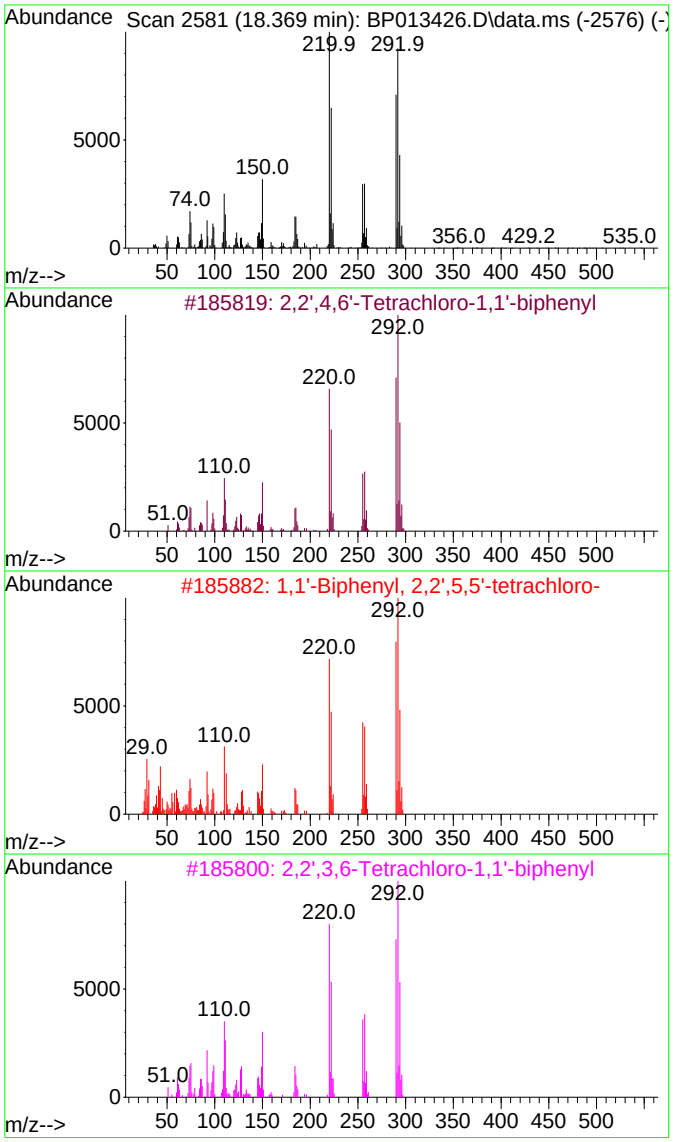
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

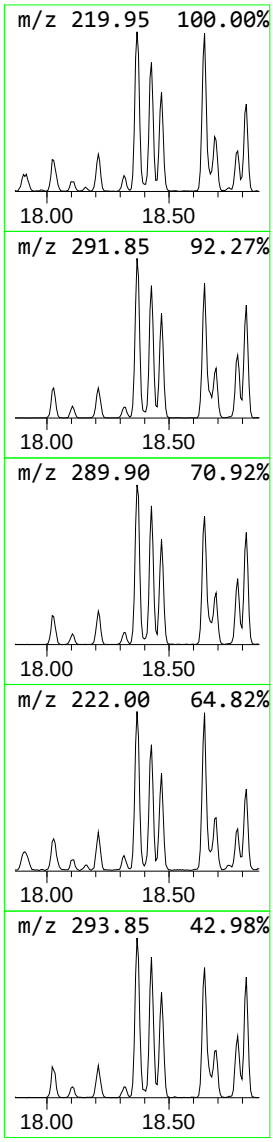
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2,2',4,6'-Tetrachloro-1,1'-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	2.69 ng/ul	862590	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
2	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
3	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

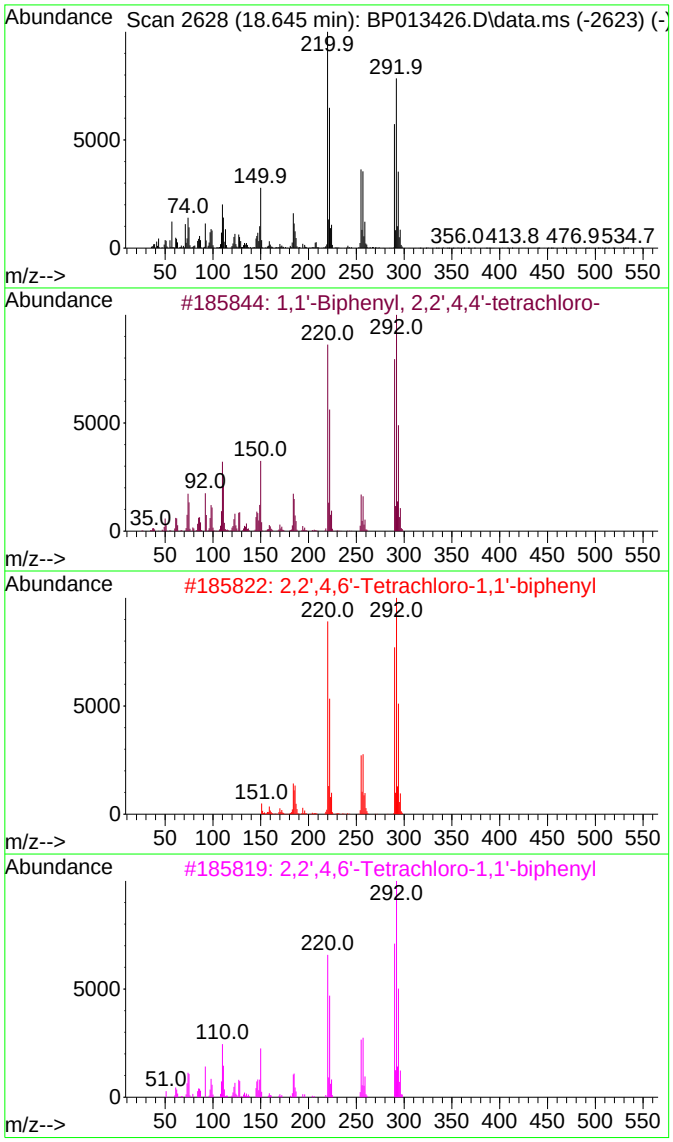
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

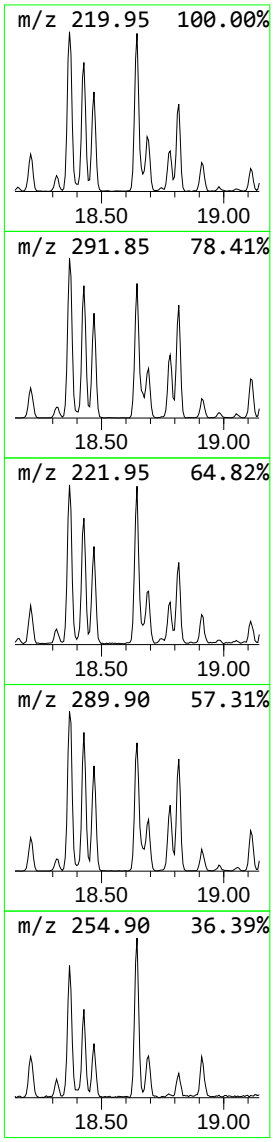
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.645	2.85 ng/ul	911742	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
3	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
5	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

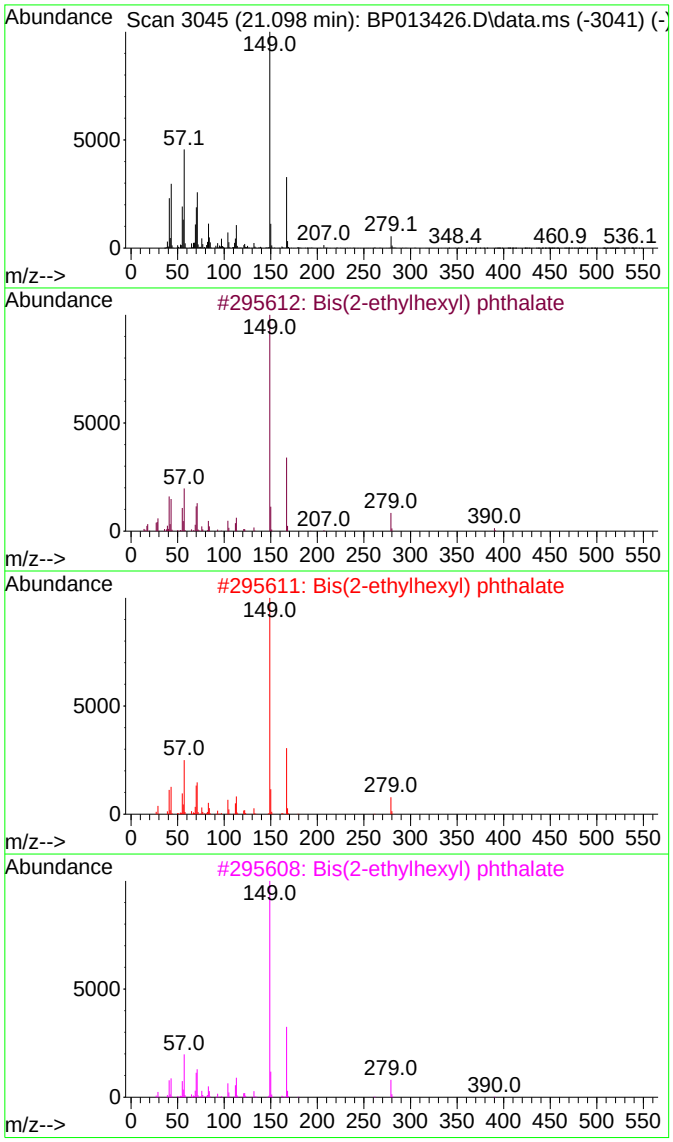
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION

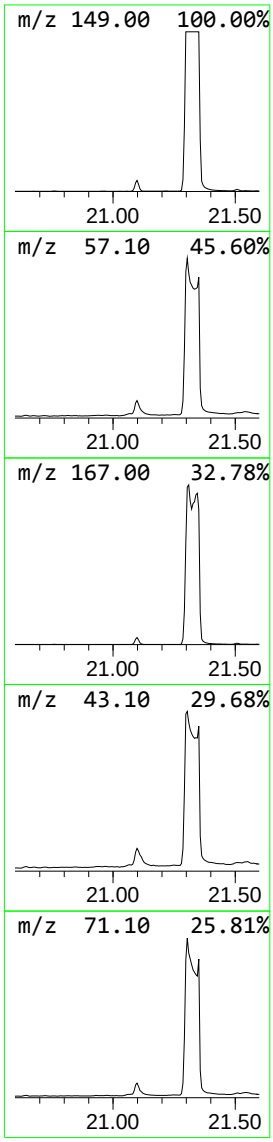
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Phthalic acid, di(6-methylh... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	11.25 ng/ul	4053890	Chrysene-d12	21.410

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	86
3	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	80
4	Phthalic acid, di(6-methylhept-2...	390	C24H38O4	1000377-97-3	72
5	Phthalic acid, di(2-propylpentyl...	390	C24H38O4	1000377-93-5	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
Data File : BP013426.D
Acq On : 10 Jan 2023 14:08
Operator : CG/JU
Sample : 01050-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4477

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.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
Cyclotrisiloxan...	4.522	2.0	ng/ul	266946	1	7.916	2656040	20.0
1,1'-Biphenyl, ...	15.839	112.0	ng/ul	29126900	3	14.563	5203720	20.0
Benzene, 1,2-di...	16.151	61.9	ng/ul	19824800	4	17.322	6402280	20.0
1,1'-Biphenyl, ...	17.192	3.0	ng/ul	948078	4	17.322	6402280	20.0
1,1'-Biphenyl, ...	17.492	2.4	ng/ul	780719	4	17.322	6402280	20.0
1,1'-Biphenyl, ...	17.910	5.4	ng/ul	1740260	4	17.322	6402280	20.0
1,1'-Biphenyl, ...	18.039	2.0	ng/ul	647341	4	17.322	6402280	20.0
n-Hexadecanoic ...	18.192	3.0	ng/ul	960910	4	17.322	6402280	20.0
2,2',4,6'-Tetra...	18.369	2.7	ng/ul	862590	4	17.322	6402280	20.0
1,1'-Biphenyl, ...	18.645	2.9	ng/ul	911742	4	17.322	6402280	20.0
Phthalic acid, ...	21.098	11.3	ng/ul	4053890	5	21.410	7206300	20.0