LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

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\SOM-EPA-BM102418MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.828	647	653	667	rVB	47957	111385	0.23%	0.119%
2	6.987	674	680	687	rBV	54940	94017	0.19%	0.100%
3	7.175	707	712	722	rVB2	66319	114453	0.23%	0.122%
4	7.640	783	791	801	rBV	773884	1266522	2.57%	1.353%
5	8.357	907	913	925	rBV	60015	127115	0.26%	0.136%
6	8.798	982	988	994	rBV2	61009	112998	0.23%	0.121%
7	9.516	1102	1110	1119	rBV2	45604	91094	0.18%	0.097%
8	10.069	1197	1204	1217	rBV5	28628	95872	0.19%	0.102%
9	10.410	1254	1262	1273	rBV	1180020	1992926	4.04%	2.129%
10	10.587	1286	1292	1301	rBV3	39167	108156	0.22%	0.116%
11	13.704	1816	1822	1829	rBV	164457	287933	0.58%	0.308%
12	13.975	1862	1868	1877	rBV	190247	326668	0.66%	0.349%
13	14.280	1912	1920	1939	rBV2	1662518	2901390	5.88%	3.100%
14	15.286	2085	2091	2097	rBV	270648	445495	0.90%	0.476%
15	15.427	2110	2115	2122	rBV5	29546	53412	0.11%	0.057%
16	15.545	2130	2135	2137	rBV	628122	919893	1.86%	0.983%
17	15.574	2137	2140	2150	rVB	1628219	2151051	4.36%	2.298%
18	15.880	2186	2192	2199	rBV	991779	1338827	2.71%	1.431%
19	15.986	2206	2210	2216	rVB	64604	95612	0.19%	0.102%
20	16.157	2234	2239	2249	rVB	151851	226559	0.46%	0.242%
21	16.269	2250	2258	2267	rBV	597688	890078	1.80%	0.951%
22	16.569	2303	2309	2315	rVB	205827	284022	0.58%	0.303%
23	16.916	2362	2368	2371	rBV	1049022	1461299	2.96%	1.561%
24	16.951	2371	2374	2379	rVV	612227	770848	1.56%	0.824%
25	17.033	2381	2388	2400	rVV2	2571870	4282693	8.68%	4.576%
26	17.133	2400	2405	2413	rVV2	353641	564684	1.14%	0.603%
27	17.210	2413	2418	2426	rVB	711480	1122110	2.27%	1.199%
28	17.492	2461	2466	2469	rBV	233310	330647	0.67%	0.353%
29	17.521	2469	2471	2478	rVB	100408	118527	0.24%	0.127%
30	17.639	2483	2491	2505	rBV	1844133	3589723	7.28%	3.836%
31	17.763	2505	2512	2519	rVV2	792687	1265919	2.57%	1.353%
32	17.833	2520	2524	2528	rVB2	60614	76689	0.16%	0.082%
33	17.886	2528	2533	2538	rBV	491206	681803	1.38%	0.728%
34	17.939	2538	2542	2549	rVB	200691	270018	0.55%	0.289%

LSC Area Percent Report

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Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

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\SOM-EPA-BM102418MA.M
Title : SVOA CALIBRATION

35	18.051	2557	2561	2565	rVB2	68568	93068	0.19%	0.099%
36	18.110	2565	2571	2577	rBV	663505	931046	1.89%	0.995%
37	18.168	2577	2581	2585	rVV	533987	698426	1.42%	0.746%
38	18.210	2585	2588	2595	rVB	457826	601529	1.22%	0.643%
39	18.392	2614	2619	2624	rBV	593316	878486	1.78%	0.939%
40	18.439	2624	2627	2633	rVV2	225780	328184	0.67%	0.351%
41	18.498	2633	2637	2640	rVV	277621	380402	0.77%	0.406%
42	18.533	2640	2643	2646	rVV	233436	324824	0.66%	0.347%
43	18.574	2646	2650	2656	rVB	460639	637439	1.29%	0.681%
44	18.674	2662	2667	2672	rVB	128336	173437	0.35%	0.185%
45	18.886	2698	2703	2707	rBV	188090	246922	0.50%	0.264%
46	18.939	2707	2712	2714	rVV	286006	426141	0.86%	0.455%
47	18.974	2714	2718	2727	rVB3	241868	410688	0.83%	0.439%
48	19.192	2750	2755	2760	rBV4	46447	91697	0.19%	0.098%
49	19.245	2760	2764	2771	rVB6	47482	81684	0.17%	0.087%
50	19.433	2791	2796	2804	rBV	438274	668676	1.36%	0.714%
51	20.951	3050	3054	3060	rBV2	196629	264312	0.54%	0.282%
52	21.162	3085	3090	3095	rBV	41304112	49329640	100.00%	52.708%
53	21.233	3097	3102	3113	rVB	3223993	4401648	8.92%	4.703%
54	23.297	3450	3453	3455	rBV	186529	263540	0.53%	0.282%
55	23.439	3471	3477	3490	rVB	1849425	3788162	7.68%	4.048%

Sum of corrected areas: 93590389

LSC Report - Integrated Chromatogram

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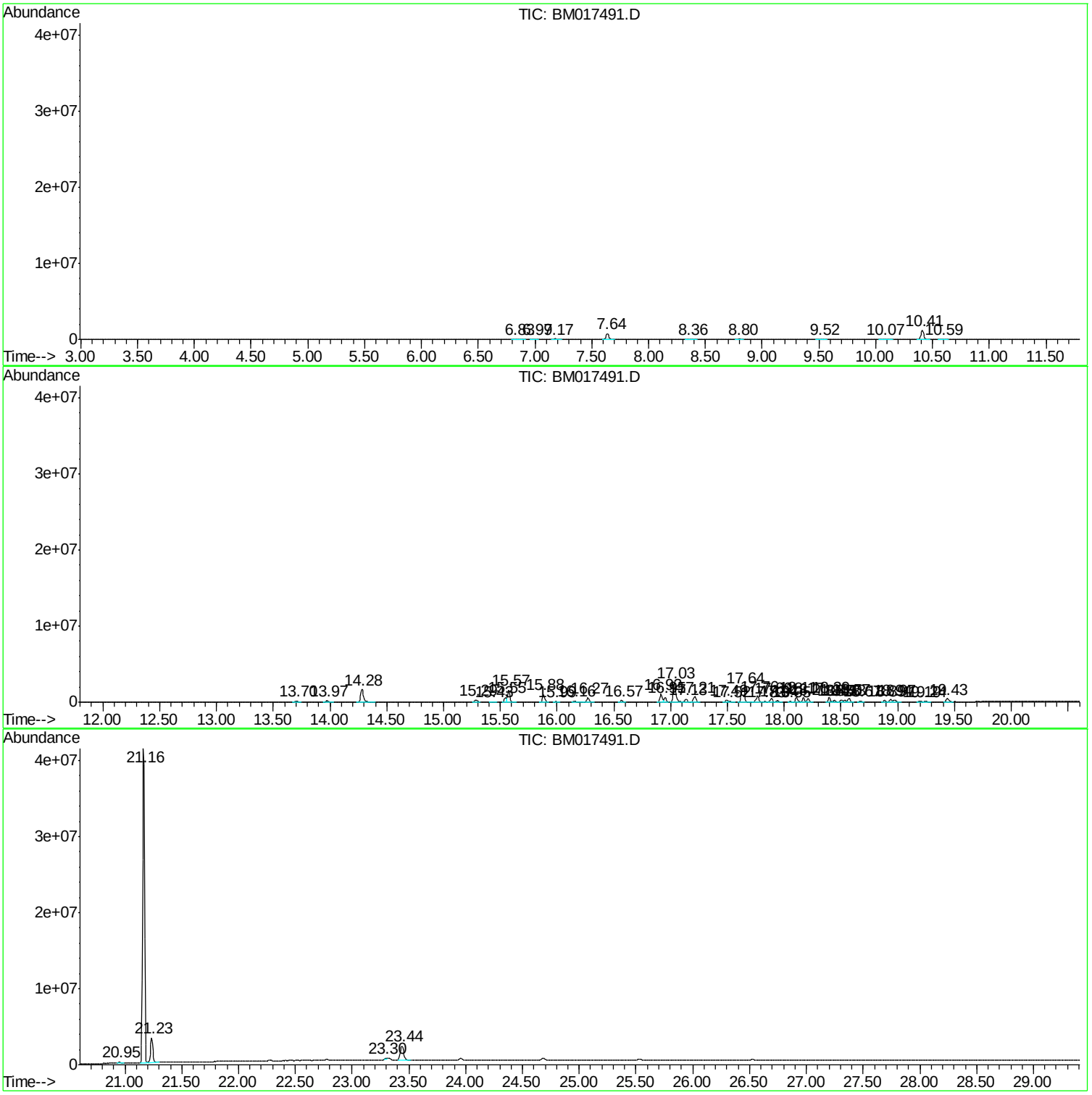
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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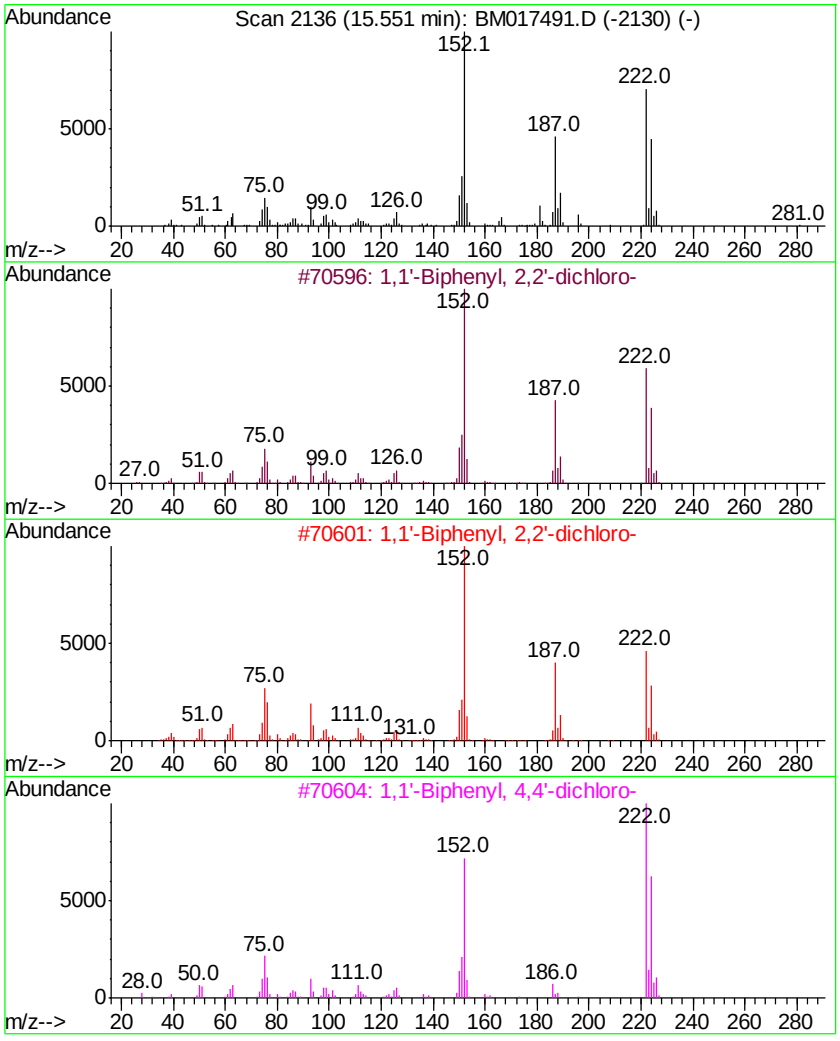
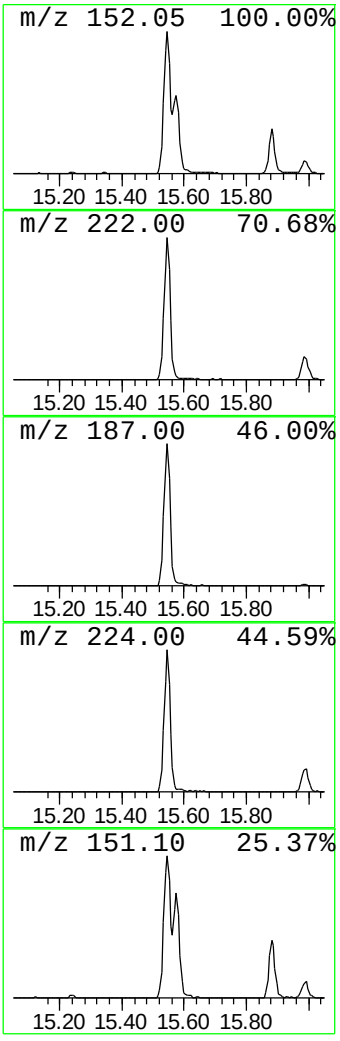
Peak Number 1 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	6.34 ng/ul	919893	Acenaphthene-d10	14.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
3	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	95
4	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	95
5	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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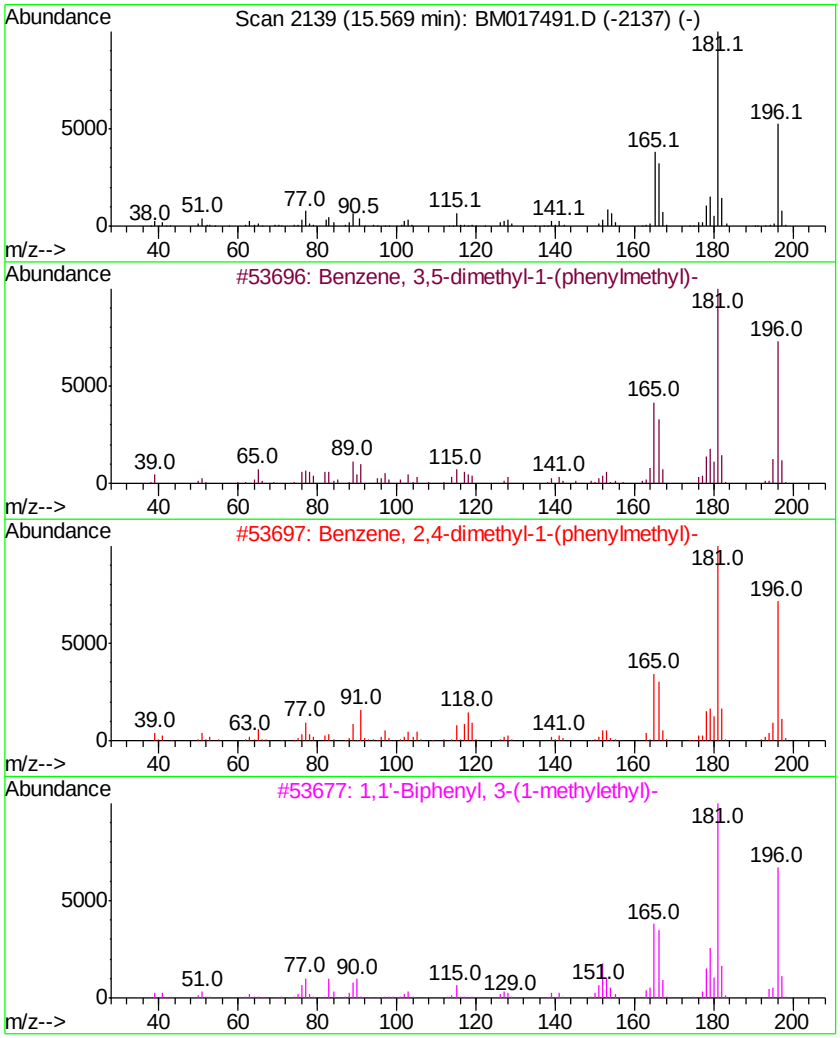
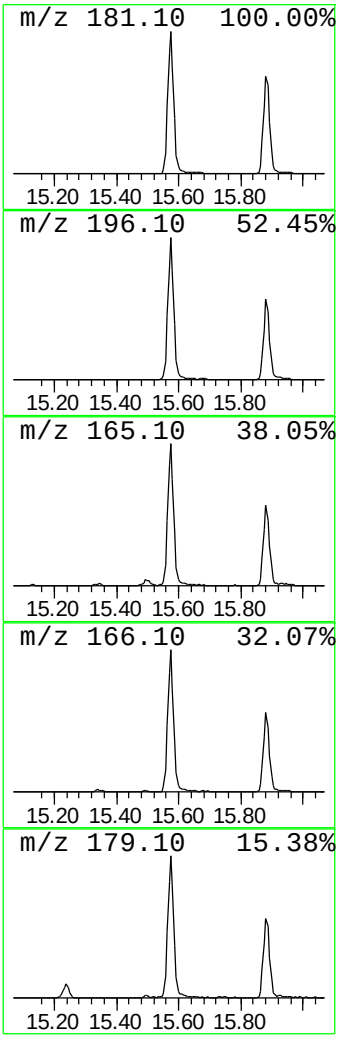
Peak Number 2 Benzene, 3,5-dimethyl-1-(ph... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	14.83 ng/ul	2151050	Acenaphthene-d10	14.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 3,5-dimethyl-1-(phenylm...	196	C15H16	028122-27-2	94
2	Benzene, 2,4-dimethyl-1-(phenylm...	196	C15H16	028122-28-3	94
3	1,1'-Biphenyl, 3-(1-methylethyl)-	196	C15H16	020282-30-8	93
4	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	93
5	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

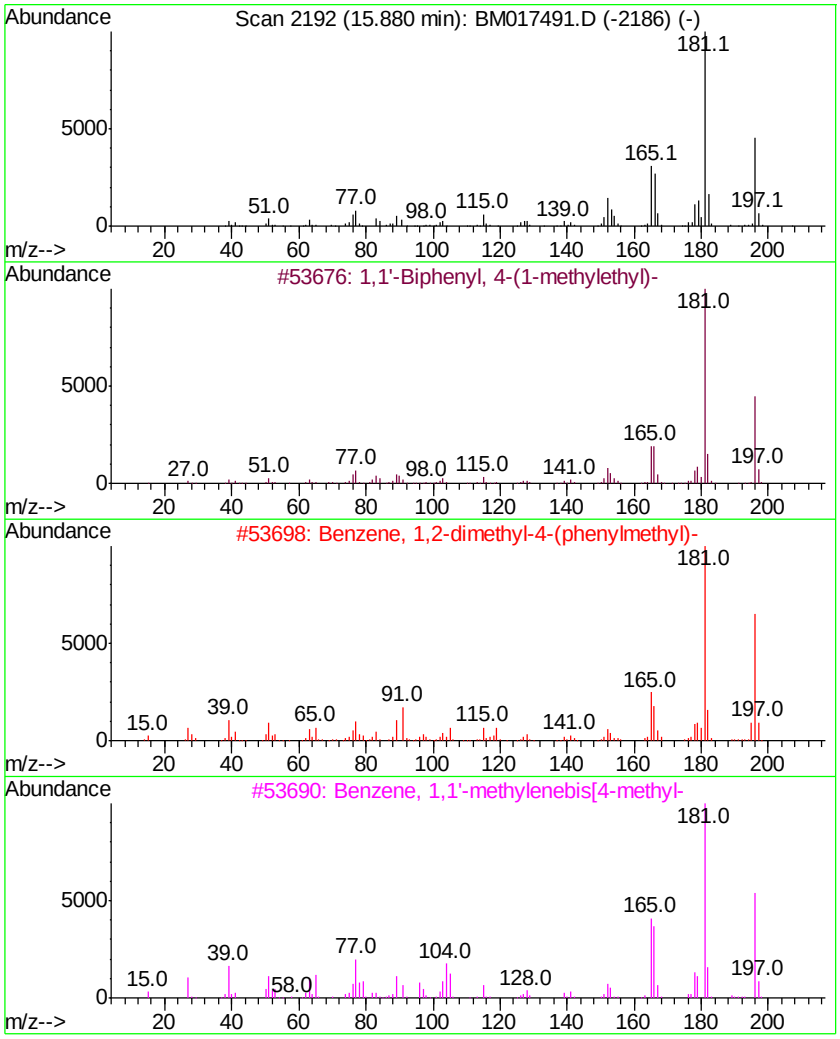
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

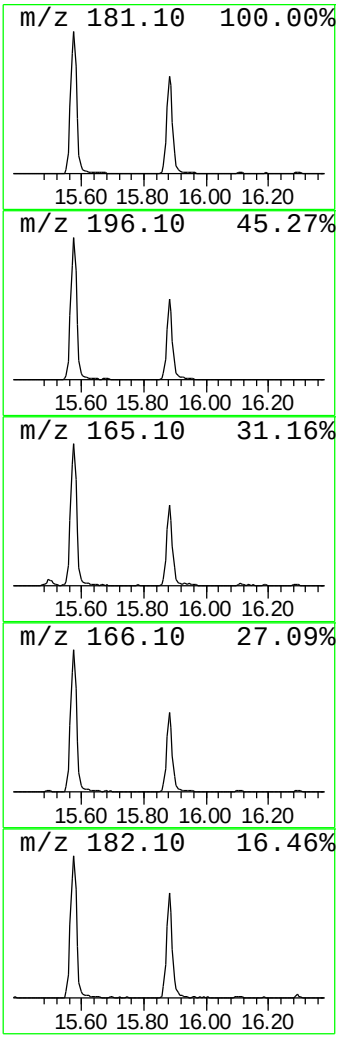
****

Peak Number 3 1,1'-Biphenyl, 4-(1-methyle... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	6.25 ng/ul	1338830	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-(1-methylethyl)-	196	C15H16	007116-95-2	96
2	Benzene, 1,2-dimethyl-4-(phenylm...	196	C15H16	013540-56-2	94
3	Benzene, 1,1'-methylenebis[4-met...	196	C15H16	004957-14-6	94
4	Benzene, 2,4-dimethyl-1-(phenylm...	196	C15H16	028122-28-3	91
5	Benzene, 1-methyl-3-[(4-methylph...	196	C15H16	021895-16-9	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
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Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

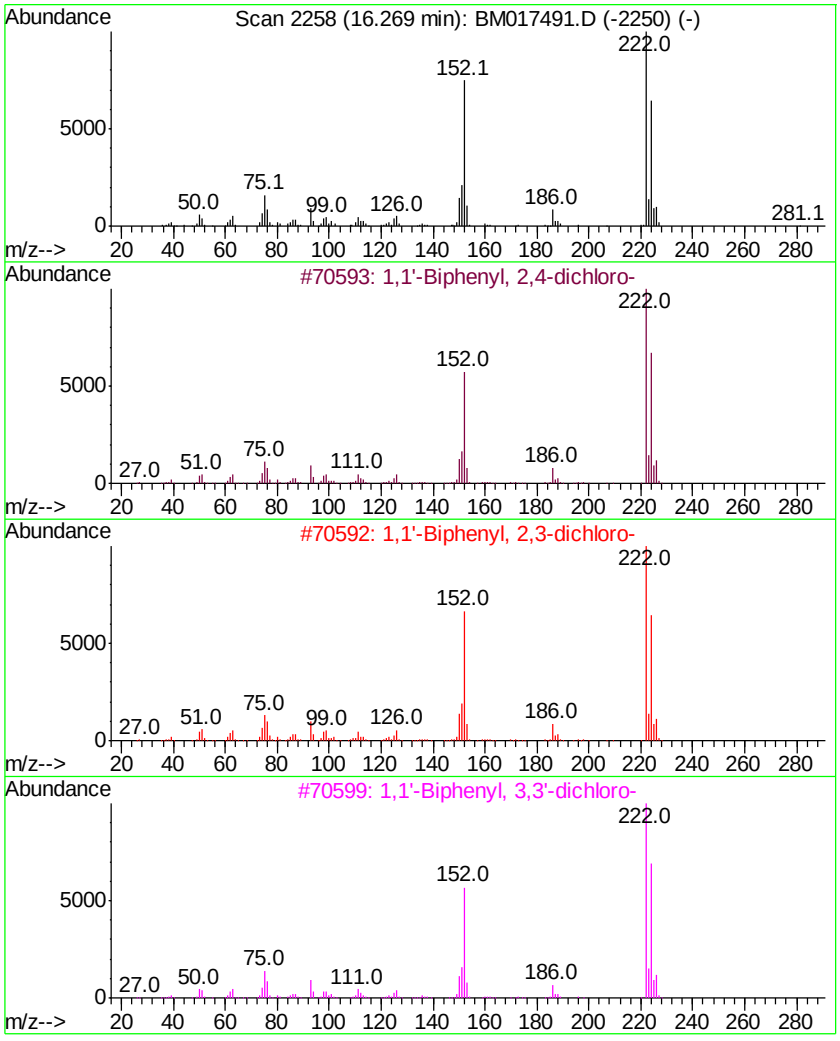
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

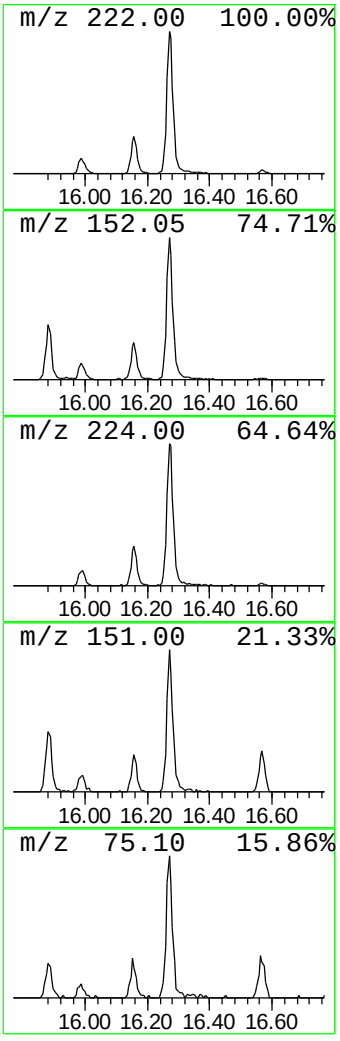
****

Peak Number 4 1,1'-Biphenyl, 2,4-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.27	4.16 ng/ul	890078	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	99
2	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99
3	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	99
4	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98
5	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	97





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ClientSampleId : A40G0MEDL

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Quant Title : SVOA CALIBRATION

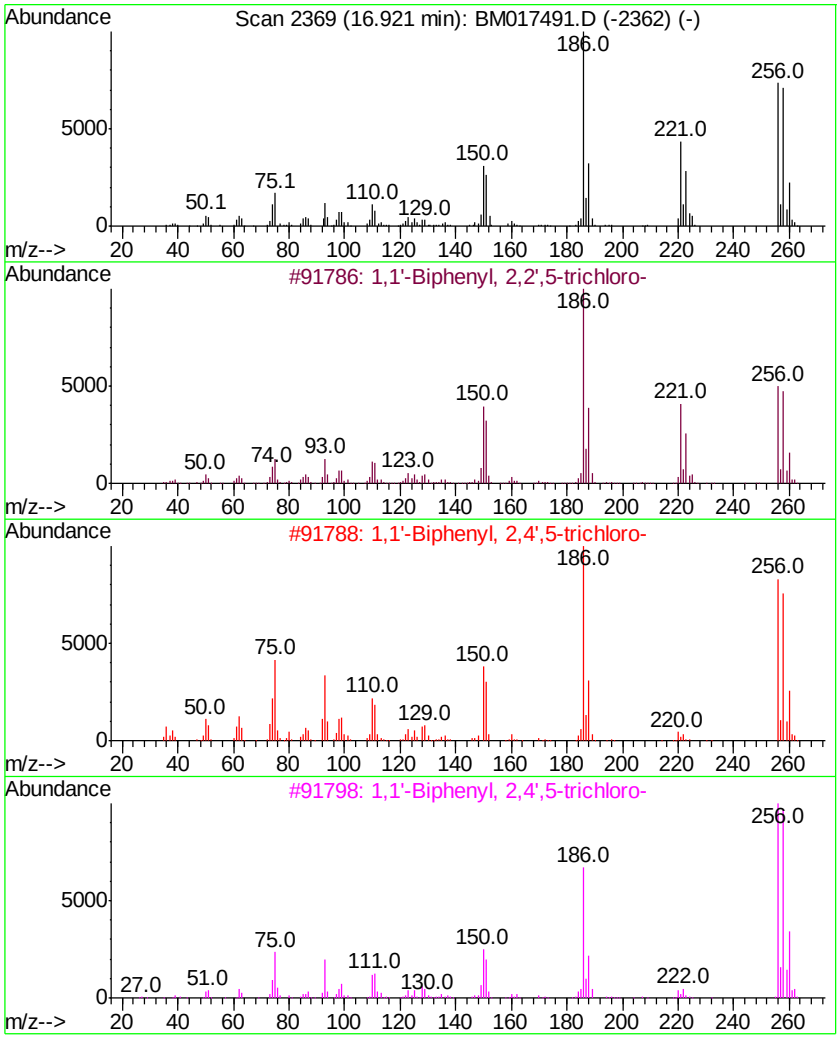
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

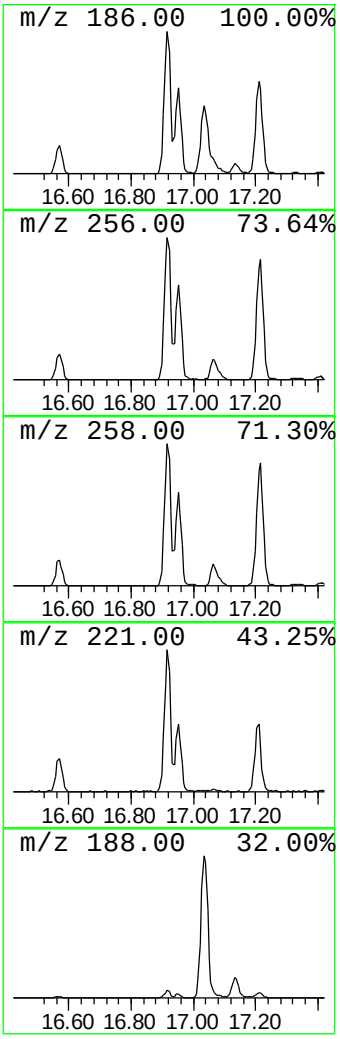
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.92	6.82 ng/ul	1461300	Phenanthrene-d10	17.03

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, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	96
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	96
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	95
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	93





Library Search Compound Report

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Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

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Quant Title : SVOA CALIBRATION

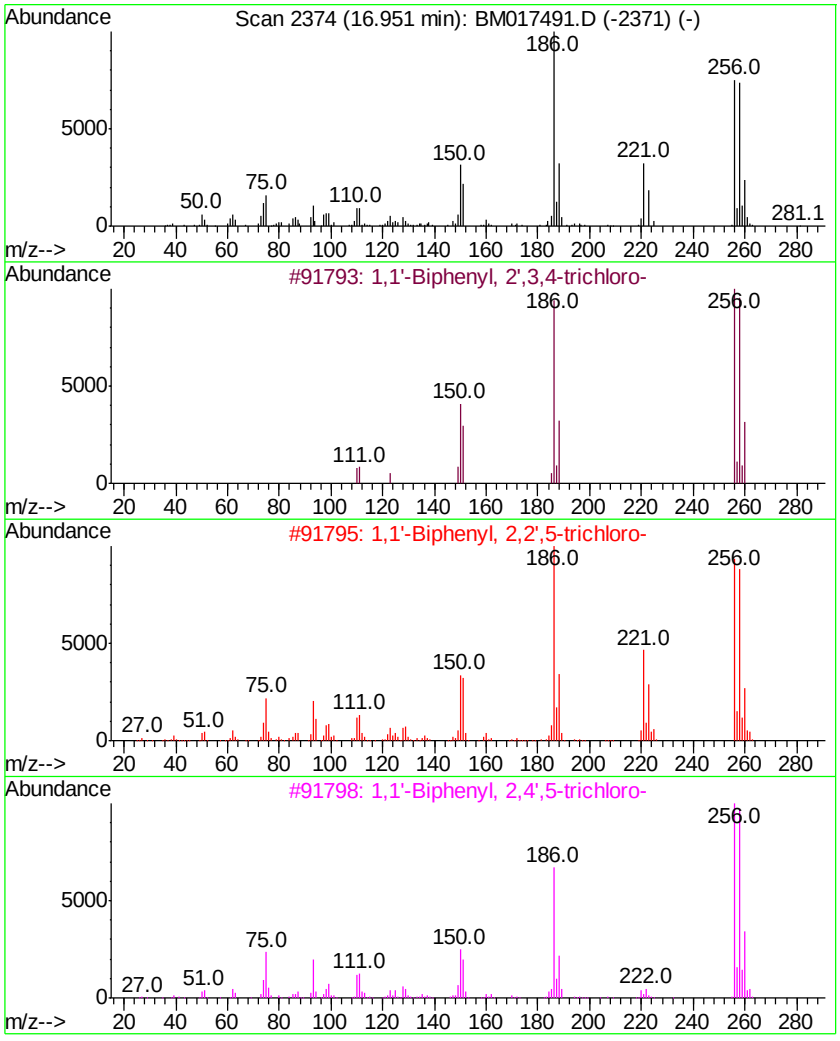
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

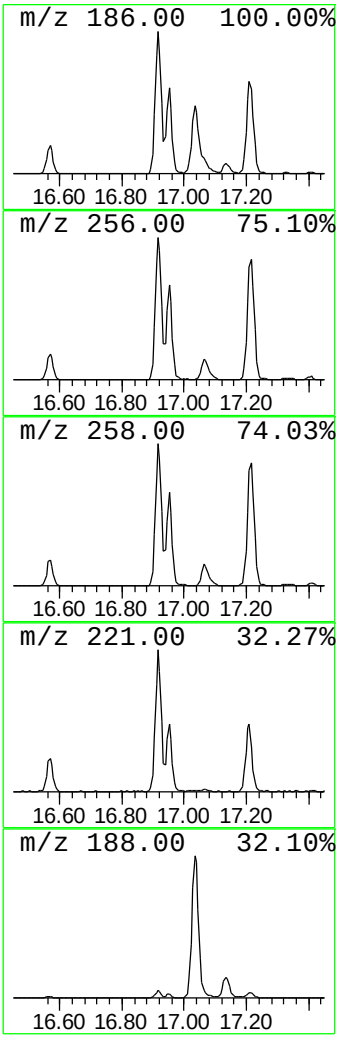
****

Peak Number 6 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.95	3.60 ng/ul	770848	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	98
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	94
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	94
4	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	94
5	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	94





Library Search Compound Report

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Operator : MJ/SJ
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

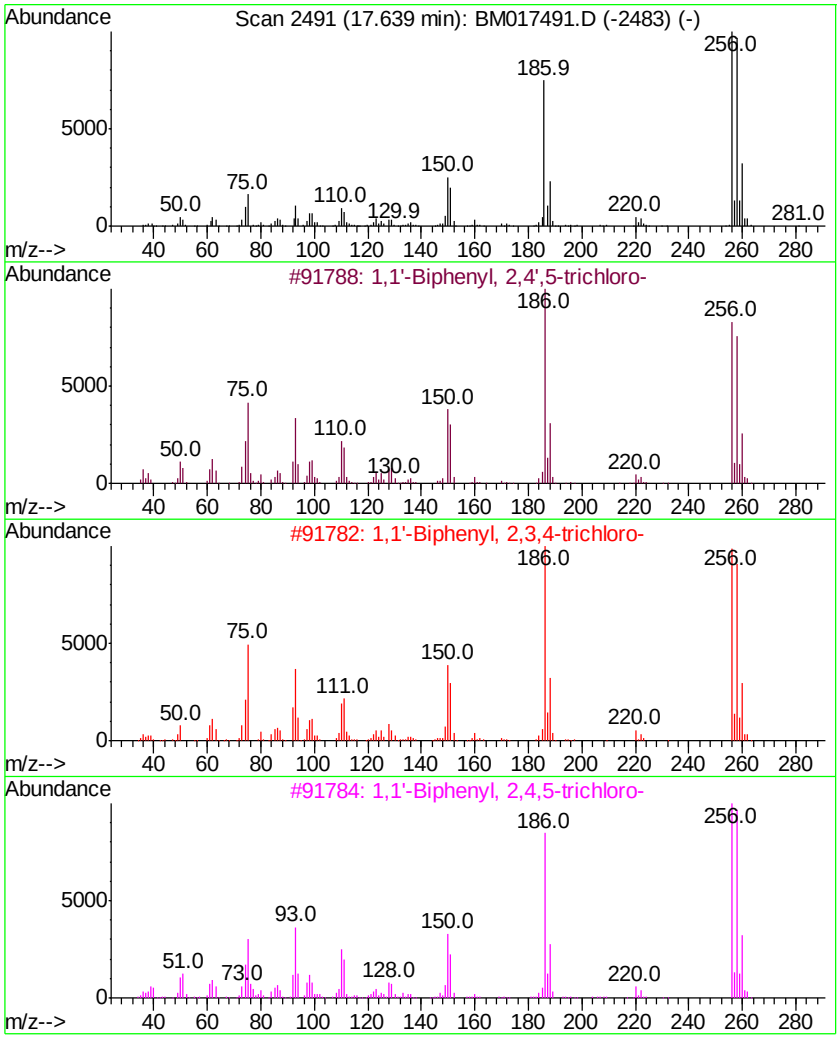
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

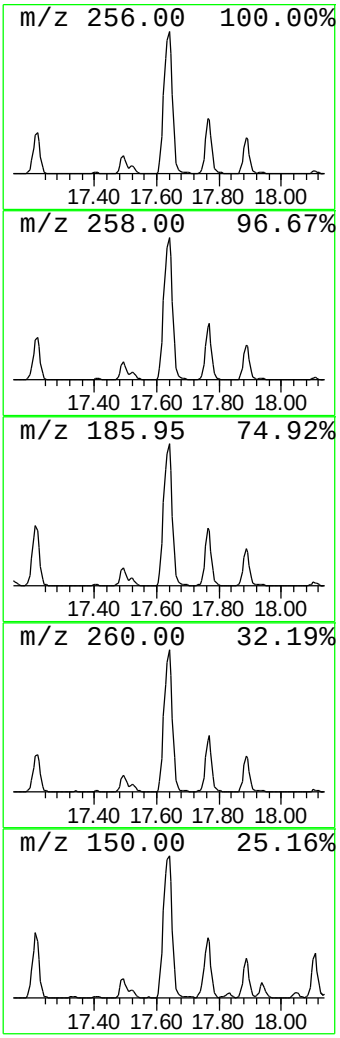
****

Peak Number 8 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.64	16.76 ng/ul	3589720	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
2	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	99
3	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98
4	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	98
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98





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Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110218\
Data File : BM017491.D
Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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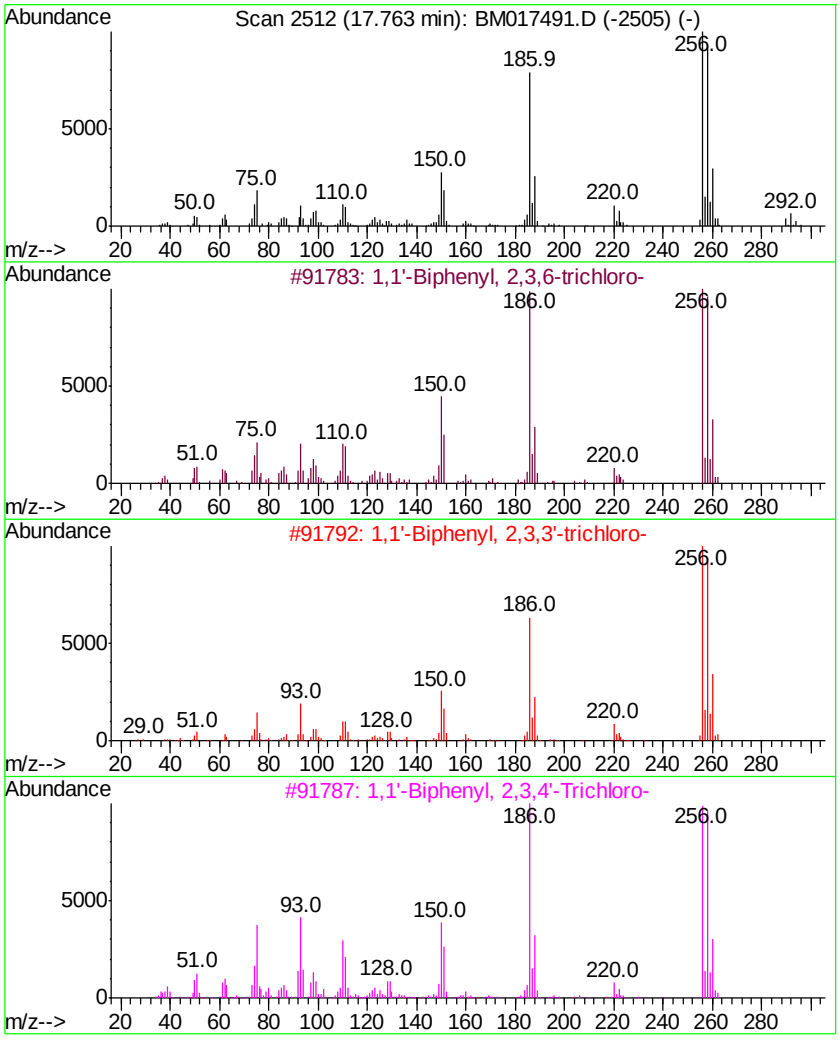
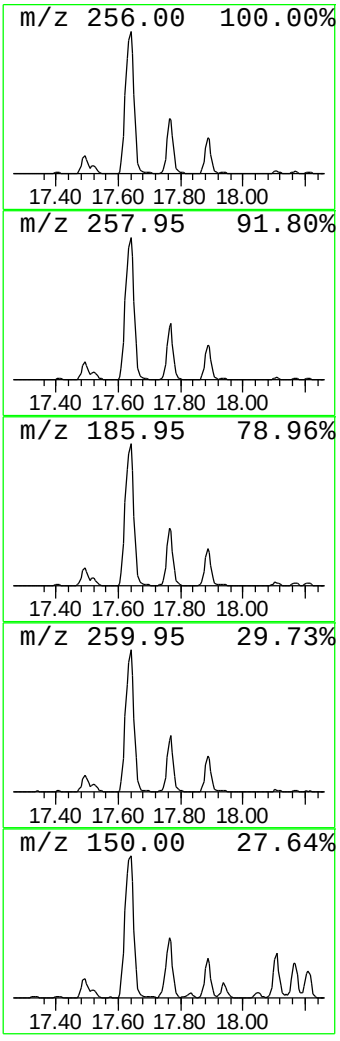
Peak Number 9 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.76	5.91 ng/ul	1265920	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	98
3	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	98
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
5	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98





Library Search Compound Report

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Data File : BM017491.D
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Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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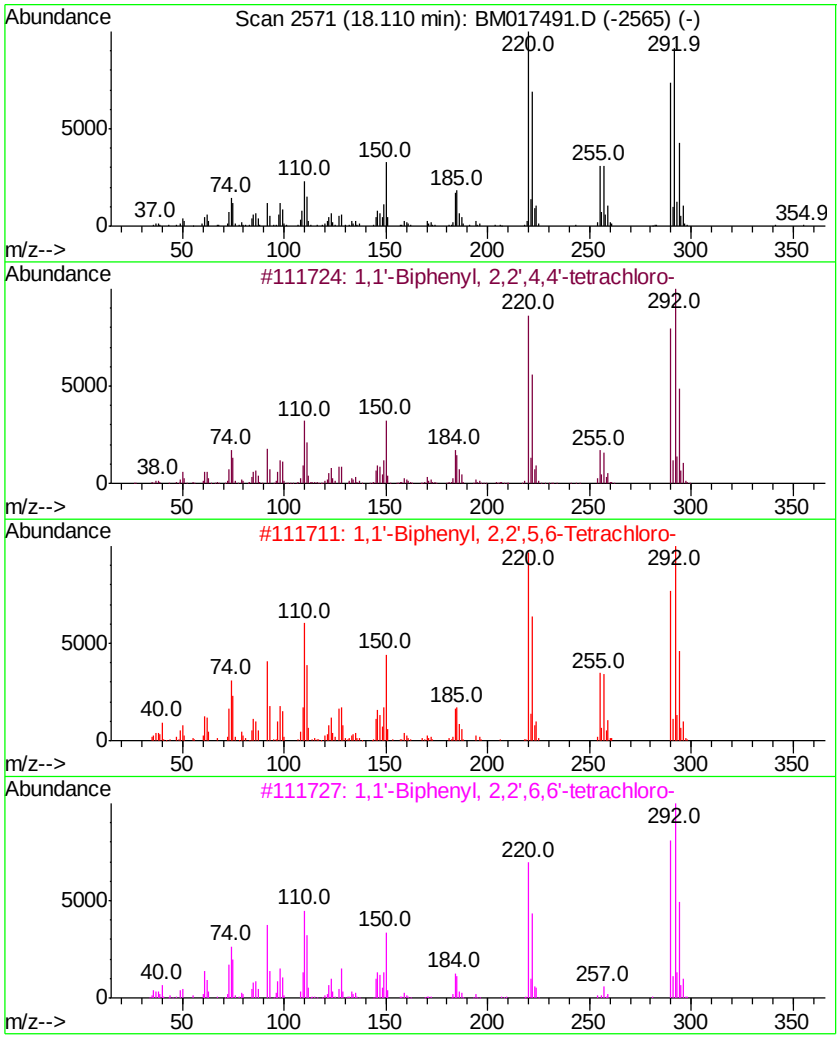
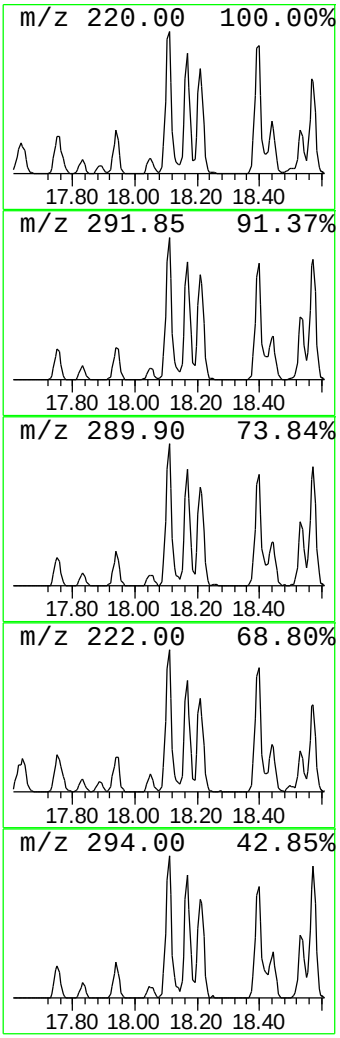
Peak Number 11 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	4.35 ng/ul	931046	Phenanthrene-d10	17.03

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	1,1'-Biphenyl, 2,2',5,6-Tetrachl...	290	C12H6Cl4	041464-41-9	99
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
4	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
5	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99





Library Search Compound Report

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Data File : BM017491.D
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Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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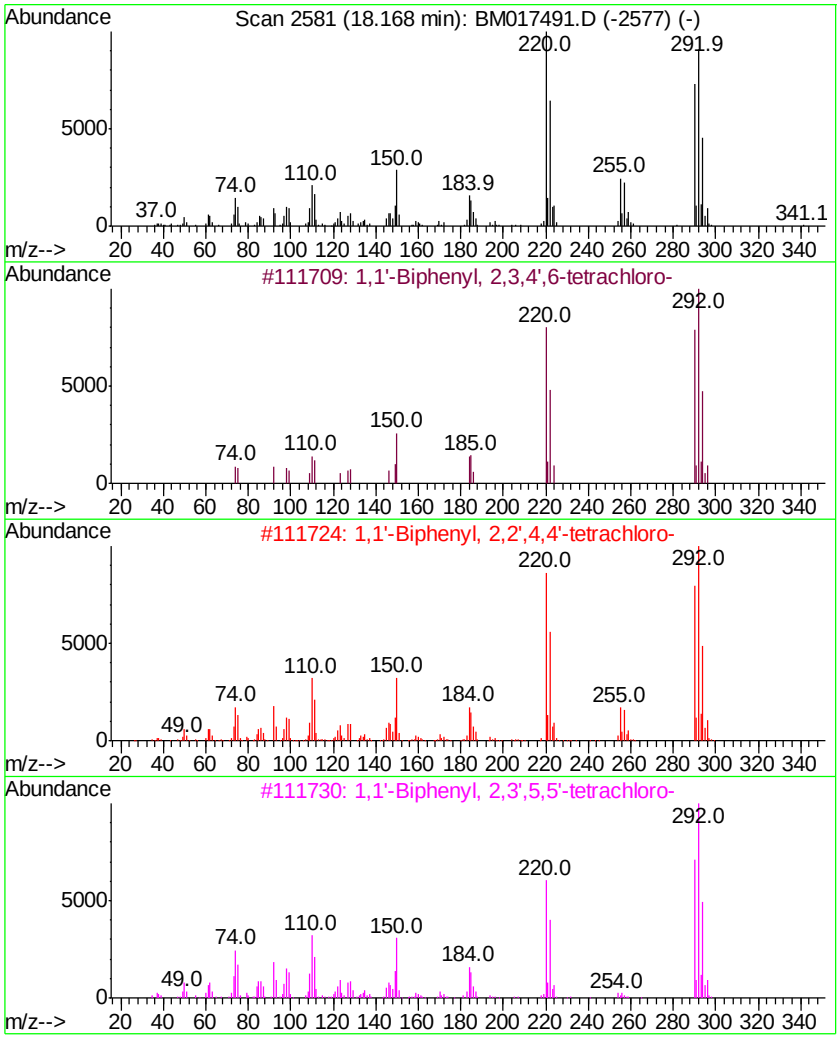
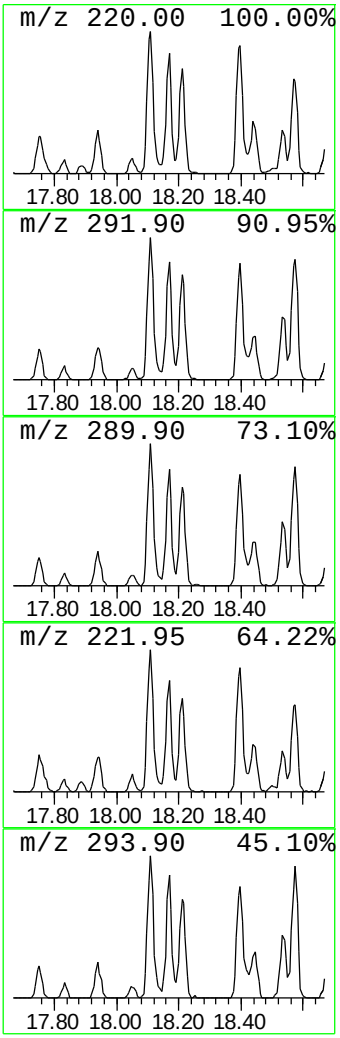
Peak Number 12 1,1'-Biphenyl, 2,3,4',6-tet... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	3.26 ng/ul	698426	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99
4	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99





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Acq On : 02 Nov 2018 11:29
Operator : MJ/SJ
Sample : J5701-16MEDL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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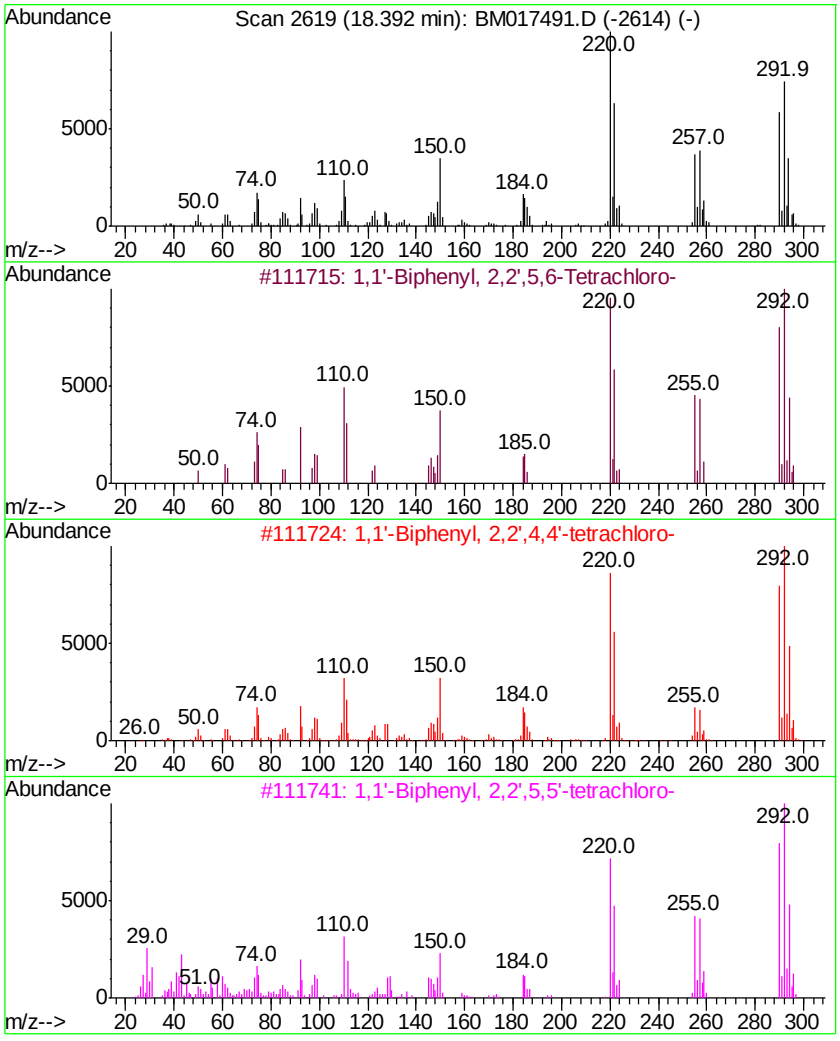
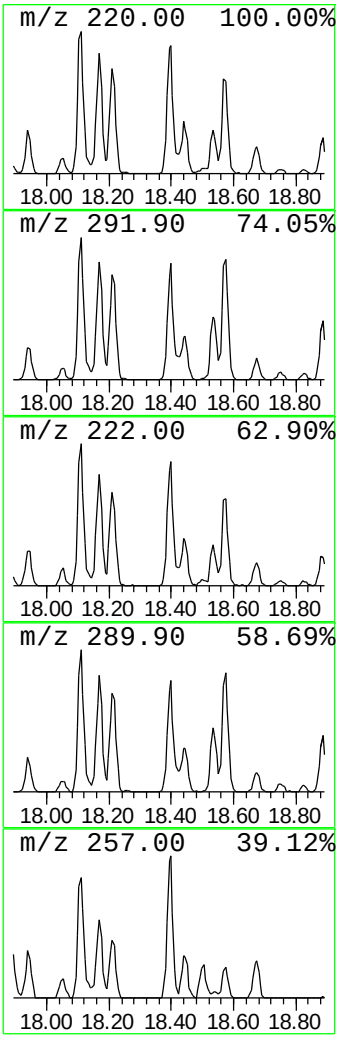
Peak Number 14 1,1'-Biphenyl, 2,2',5,6-Tet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.39	4.10 ng/ul	878486	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,6-Tetrachl...	290	C12H6Cl4	041464-41-9	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
4	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
5	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110218\
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Acq On : 02 Nov 2018 11:29
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Instrument : BNA_M
ClientSampleId : A40G0MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1,1'-Biphenyl, 2,...	15.55	6.3	ng/ul	919893	3	14.28	2901390	20.0
Benzene, 3,5-dime...	15.57	14.8	ng/ul	2151050	3	14.28	2901390	20.0
1,1'-Biphenyl, 4-...	15.88	6.3	ng/ul	1338830	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	16.27	4.2	ng/ul	890078	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	16.92	6.8	ng/ul	1461300	4	17.03	4282690	20.0
1,1'-Biphenyl, 2'...	16.95	3.6	ng/ul	770848	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	17.64	16.8	ng/ul	3589720	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	17.76	5.9	ng/ul	1265920	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	18.11	4.3	ng/ul	931046	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	18.17	3.3	ng/ul	698426	4	17.03	4282690	20.0
1,1'-Biphenyl, 2,...	18.39	4.1	ng/ul	878486	4	17.03	4282690	20.0