LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.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	9.934	1173	1181	1198	rBV	698149	1513908	1.05%	0.499%
2	13.592	1798	1803	1808	rBV	1504903	2059345	1.42%	0.679%
3	13.863	1839	1849	1860	rBV	1836674	3056478	2.11%	1.007%
4	14.175	1893	1902	1909	rBV2	1289422	2245701	1.55%	0.740%
5	15.175	2067	2072	2077	rBV	2187111	3081031	2.13%	1.015%
6	16.933	2366	2371	2375	rBV	1586203	2375205	1.64%	0.783%
7	16.980	2375	2379	2383	rVB	3097279	3967034	2.74%	1.307%
8	17.033	2383	2388	2392	rBV	2574146	3463539	2.39%	1.141%
9	17.857	2525	2528	2534	rVB	1224501	1658619	1.15%	0.547%
10	18.375	2612	2616	2622	rVB	1421060	1919402	1.33%	0.632%
11	19.010	2720	2724	2731	rBV	4723693	6609990	4.57%	2.178%
12	19.151	2745	2748	2753	rVB4	1098393	1481172	1.02%	0.488%
13	19.351	2777	2782	2784	rBV2	3664803	5176301	3.58%	1.706%
14	19.380	2784	2787	2791	rVB	4450097	5216557	3.61%	1.719%
15	19.492	2802	2806	2810	rVB	1507791	1804159	1.25%	0.594%
16	20.010	2890	2894	2898	rVV	10764432	11963165	8.27%	3.942%
17	20.204	2923	2927	2929	rBV	2016859	2465001	1.70%	0.812%
18	20.233	2929	2932	2939	rVB2	3382754	4221430	2.92%	1.391%
19	20.369	2952	2955	2959	rVB	1600956	1881664	1.30%	0.620%
20	20.451	2964	2969	2977	rVB	18094405	20555713	14.21%	6.773%
21	20.874	3038	3041	3049	rVB5	1525733	2073080	1.43%	0.683%
22	21.098	3071	3079	3082	rBV2	104807970	144670209	100.00%	47.670%
23	21.151	3082	3088	3093	rVV3	2761255	6630294	4.58%	2.185%
24	21.192	3093	3095	3100	rVB	2315970	2675811	1.85%	0.882%
25	21.374	3122	3126	3129	rBV	4254527	5423696	3.75%	1.787%
26	21.415	3129	3133	3137	rVB	21687893	25306733	17.49%	8.339%
27	21.463	3138	3141	3145	rVB2	1282492	1451524	1.00%	0.478%
28	21.627	3166	3169	3173	rBV	3928536	4553414	3.15%	1.500%
29	22.286	3278	3281	3287	rVB2	2055053	3073142	2.12%	1.013%
30	22.704	3348	3352	3362	rVB	1899202	5700330	3.94%	1.878%
31	22.998	3398	3402	3408	rVB	1900079	3152159	2.18%	1.039%
32	25.592	3838	3843	3857	rVB	2388876	6329376	4.38%	2.086%
33	26.850	4050	4057	4067	rVB2	1918770	5728380	3.96%	1.888%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Title : SVOA CALIBRATION

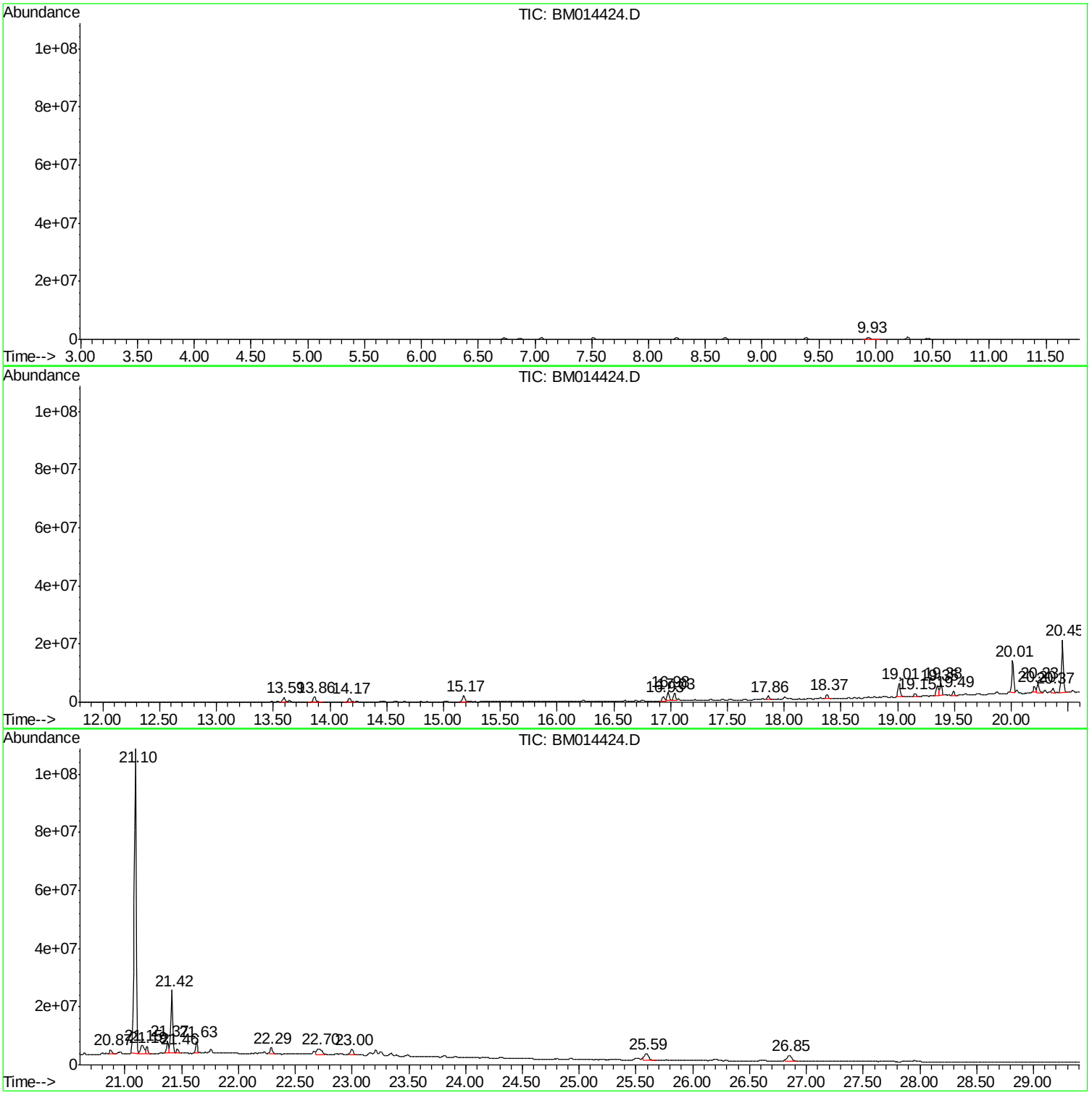
Sum of corrected areas: 303483562

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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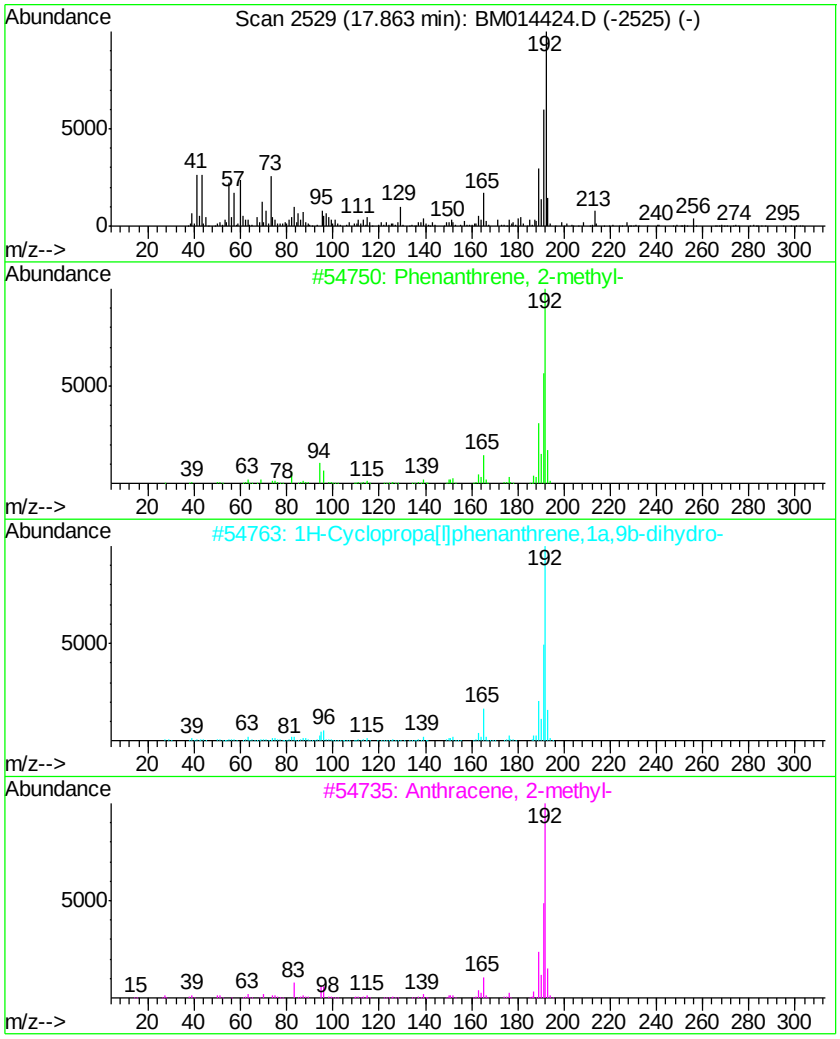
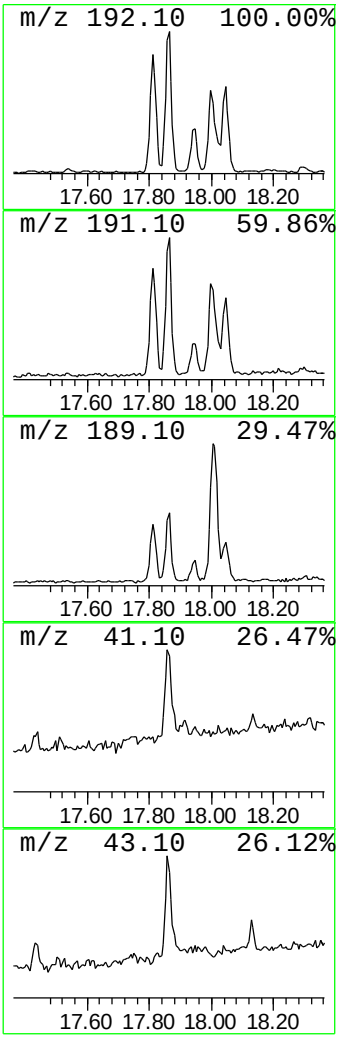
****

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.86	13.97 ng/ul	1658620	Phenanthrene-d10	16.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	92





Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

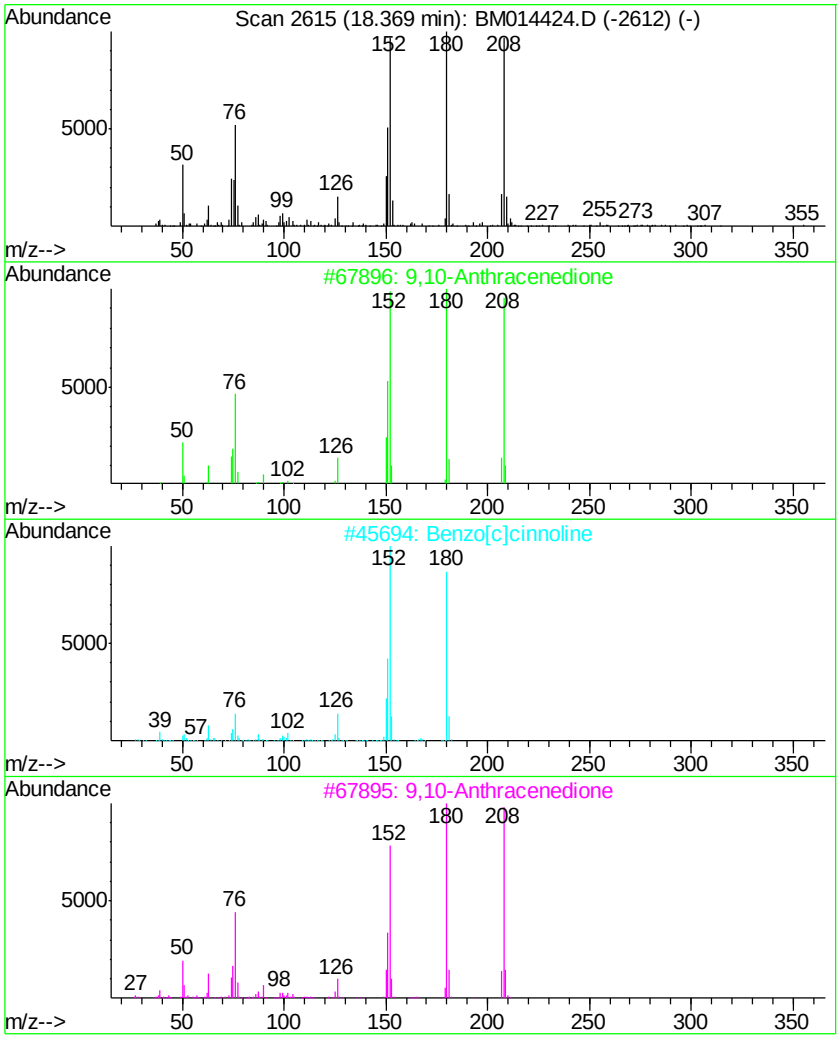
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

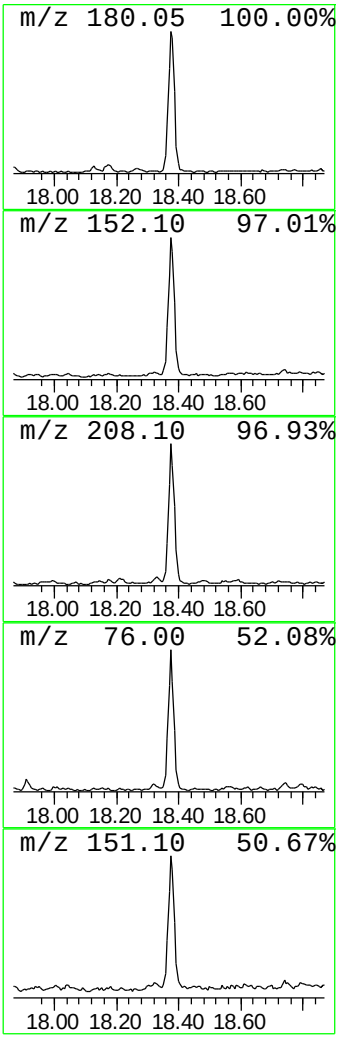
****

Peak Number 2 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	16.16 ng/ul	1919400	Phenanthrene-d10	16.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	95
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	90
5	9H-Fluoren-9-one	180	C13H8O	000486-25-9	87





Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
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Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

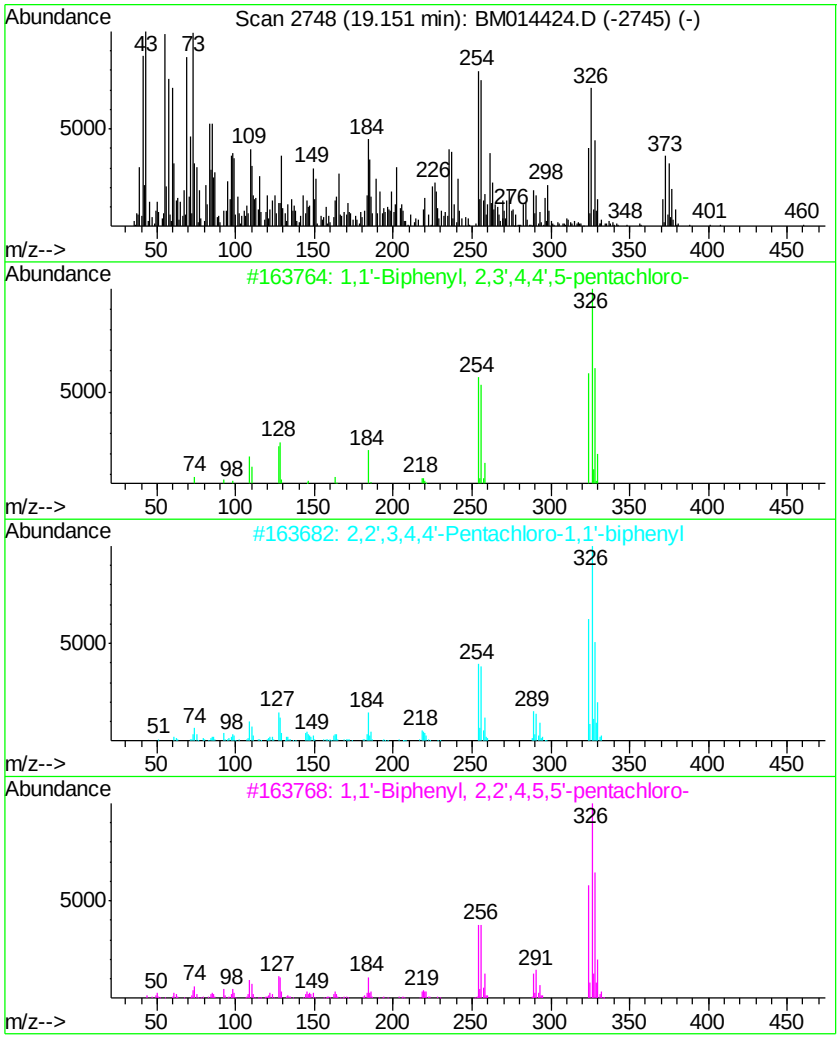
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

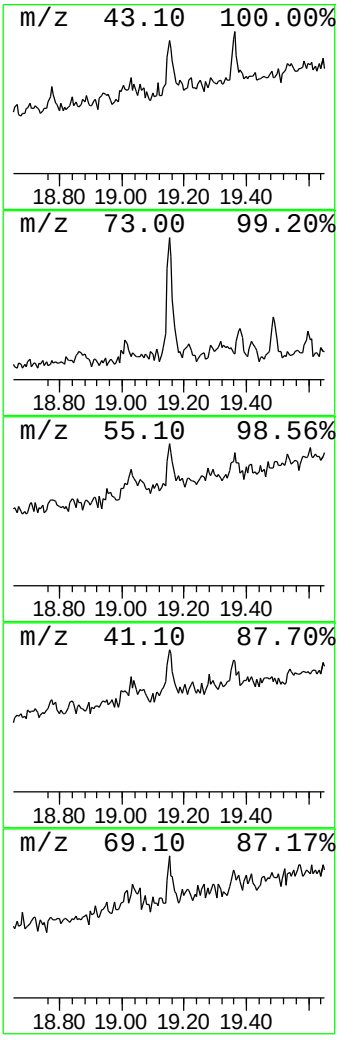
****

Peak Number 3 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.15	4.47 ng/ul	1481170	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	97
2	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	93
3	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	93
4	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	93
5	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	93





Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
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Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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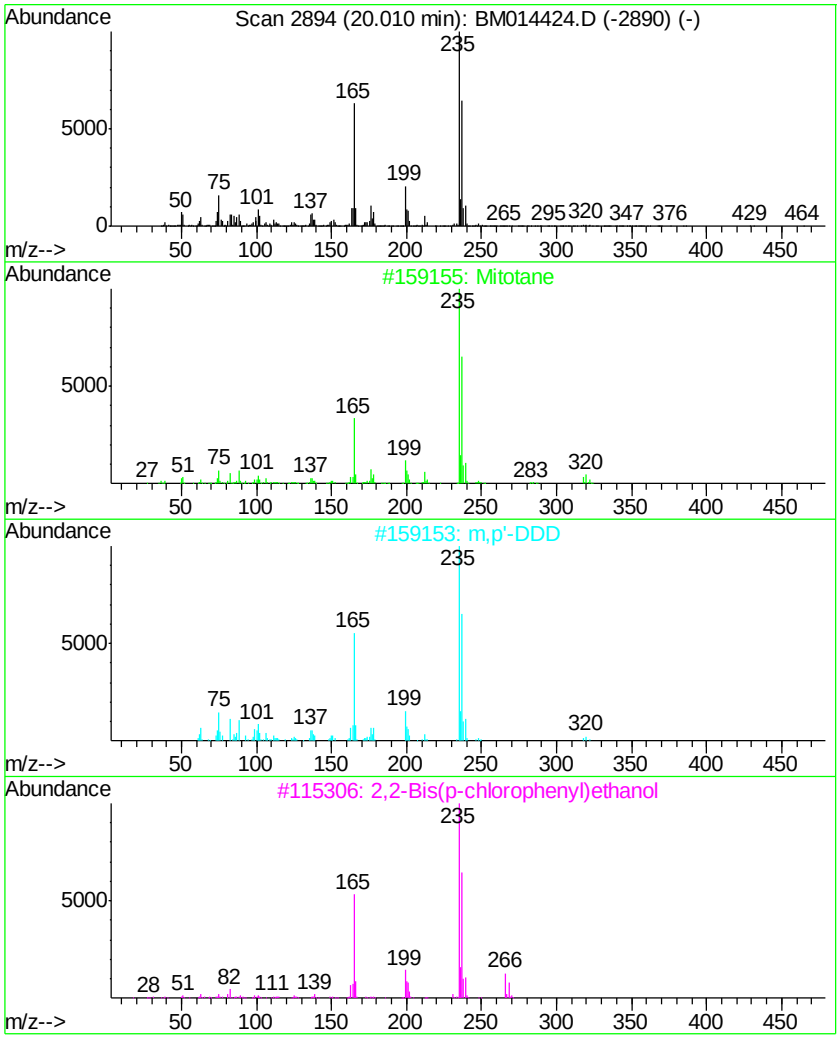
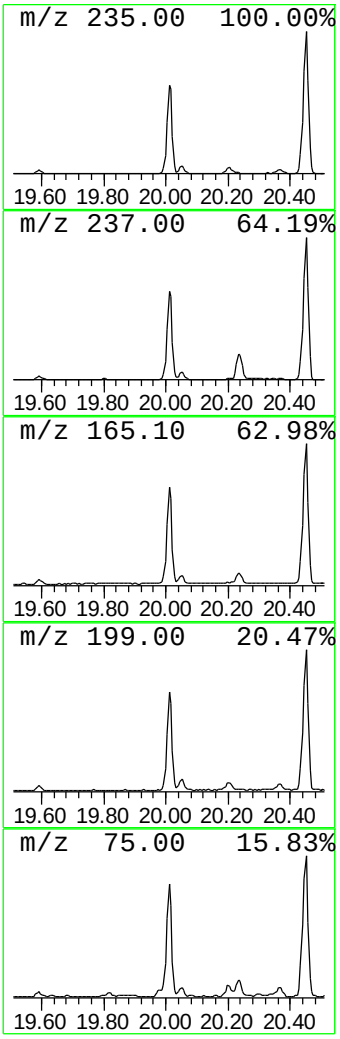
Peak Number 5 Mitotane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.01	36.09 ng/ul	11963200	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mitotane	318	C14H10Cl4	000053-19-0	99
2	m,p'-DDD	318	C14H10Cl4	004329-12-8	91
3	2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	86
4	2,2-Bis(4-chlorophenyl)acetic acid	280	C14H10Cl2O2	000083-05-6	86
5	3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	86





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Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

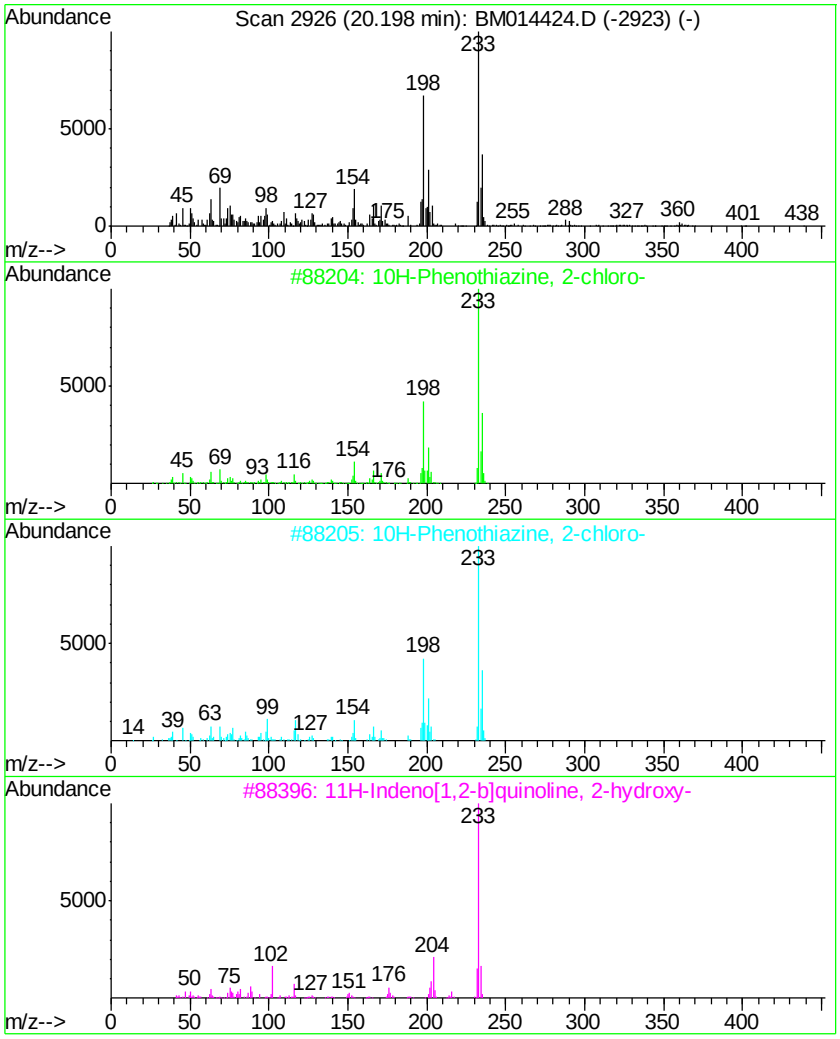
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

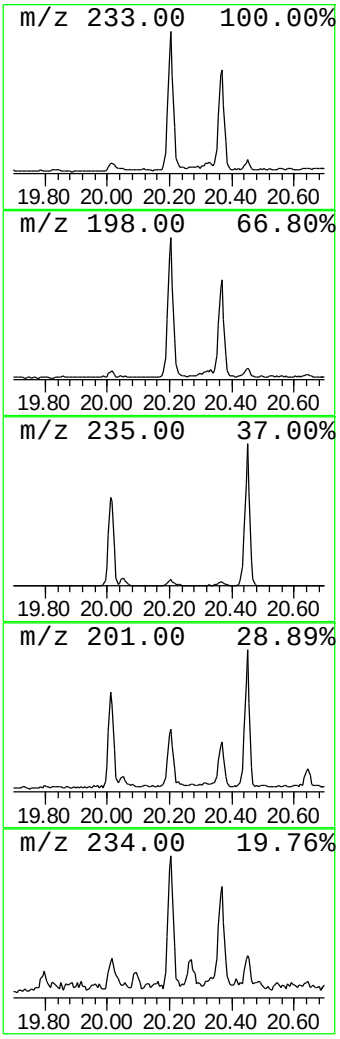
****

Peak Number 6 10H-Phenothiazine, 2-chloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.20	7.44 ng/ul	2465000	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10H-Phenothiazine, 2-chloro-	233	C12H8ClNS	000092-39-7	99
2	10H-Phenothiazine, 2-chloro-	233	C12H8ClNS	000092-39-7	93
3	11H-Indeno[1,2-b]quinoline, 2-hy...	233	C16H11NO	1000214-19-5	46
4	1H-Phenanthro[9,10-d]imidazol-2-...	233	C15H11N3	037052-13-4	38
5	Phenanthro[9,10-d]oxazole, 2-met...	233	C16H11NO	021639-88-3	38





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Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

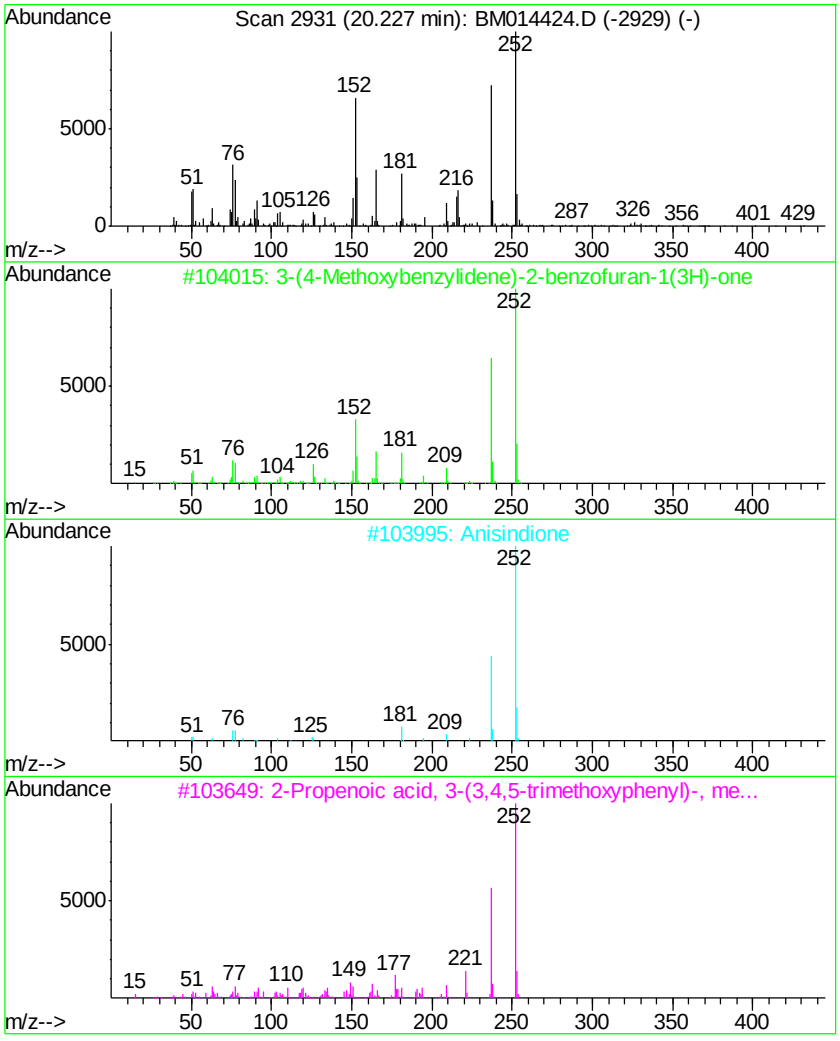
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

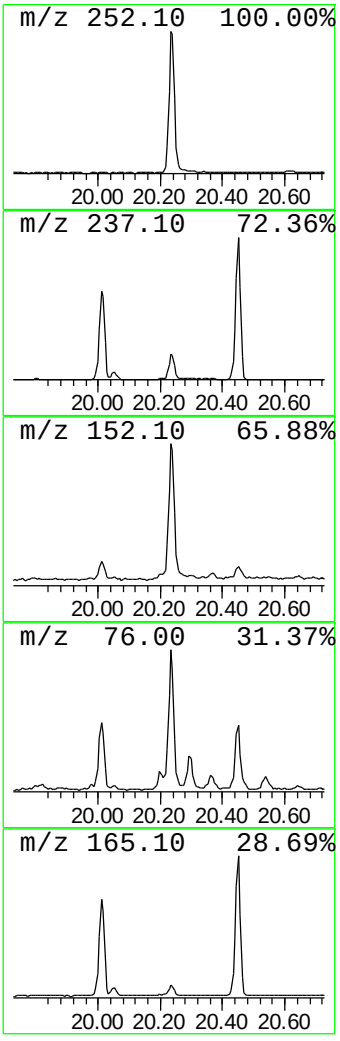
****

Peak Number 7 3-(4-Methoxybenzylidene)-2-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.23	12.73 ng/ul	4221430	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(4-Methoxybenzylidene)-2-benzo...	252	C16H12O3	1000332-19-6	96
2	Anisindione	252	C16H12O3	000117-37-3	83
3	2-Propenoic acid, 3-(3,4,5-trime...	252	C13H16O5	007560-49-8	43
4	3H-Pyrrolo[3,2-f]quinoline, 1,2,...	252	C17H20N2	1000338-03-5	43
5	2-Propenoic acid, 3-(3,4,5-trime...	252	C13H16O5	007560-49-8	38





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Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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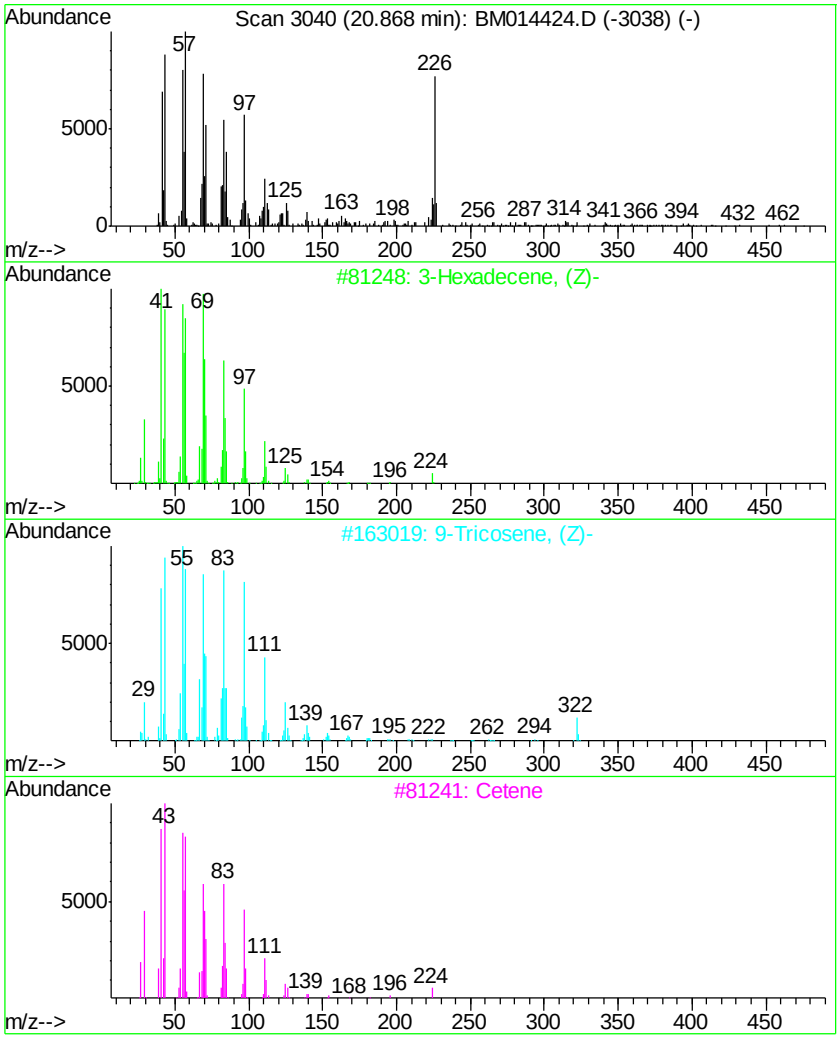
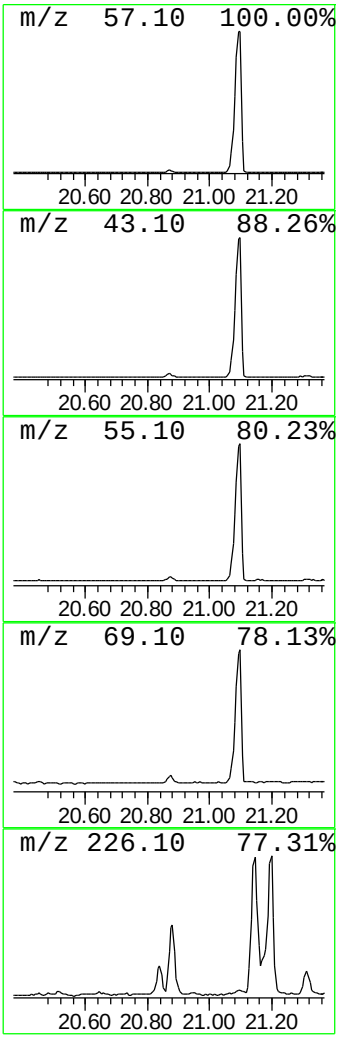
Peak Number 10 (DEL) Alkane: Cyclic20.87 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	6.25 ng/ul	2073080	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexadecene, (Z)-	224	C16H32	034303-81-6	70
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	62
3	Cetene	224	C16H32	000629-73-2	49
4	Octacosanol	410	C28H58O	000557-61-9	46
5	1-Octadecene	252	C18H36	000112-88-9	46





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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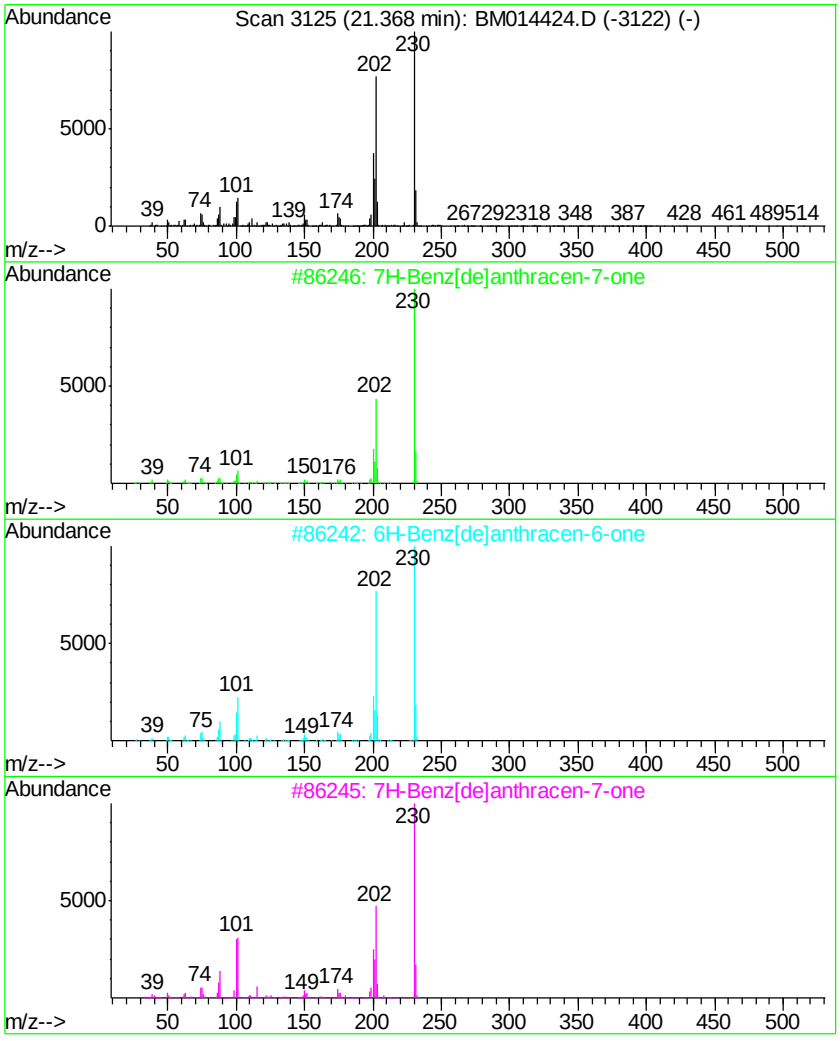
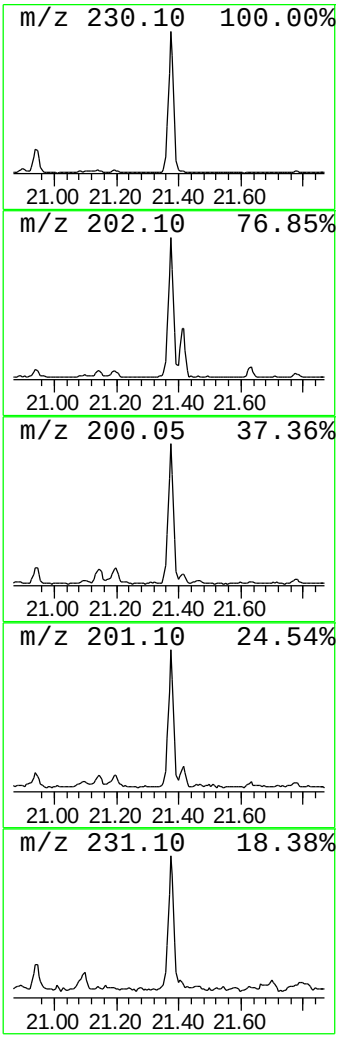
Peak Number 11 7H-Benz[de]anthracen-7-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	16.36 ng/ul	5423700	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	98
2	6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	97
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	97
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	96
5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	94





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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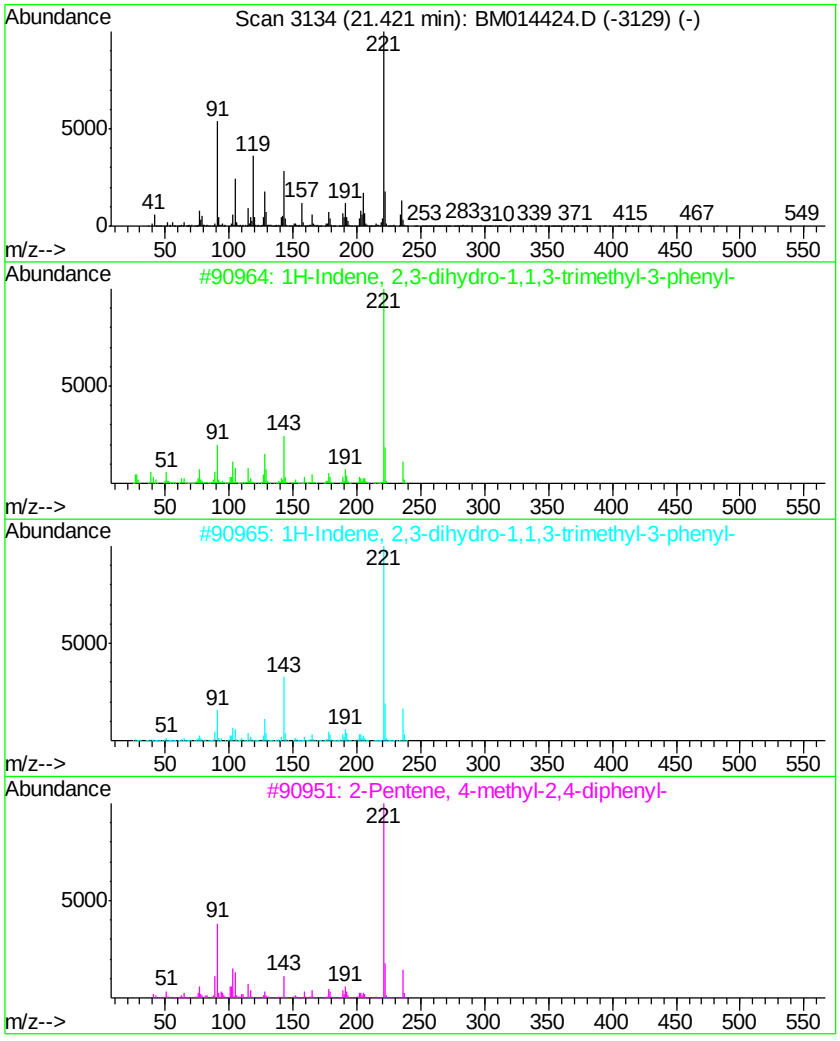
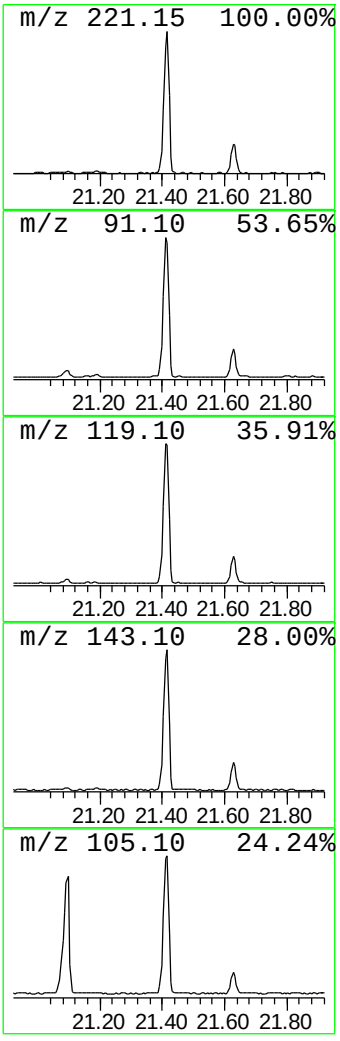
Peak Number 12 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.42	76.34 ng/ul	25306700	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	52
2	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	47
3	2-Pentene, 4-methyl-2,4-diphenyl-	236	C18H20	006258-73-7	37
4	4-Hydroxy-beta-phenylcinnamonitrile	221	C15H11NO	016281-90-6	35
5	Methanone, cyclohexyl(2,5-diprop...	304	C19H28O3	076185-90-5	35





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Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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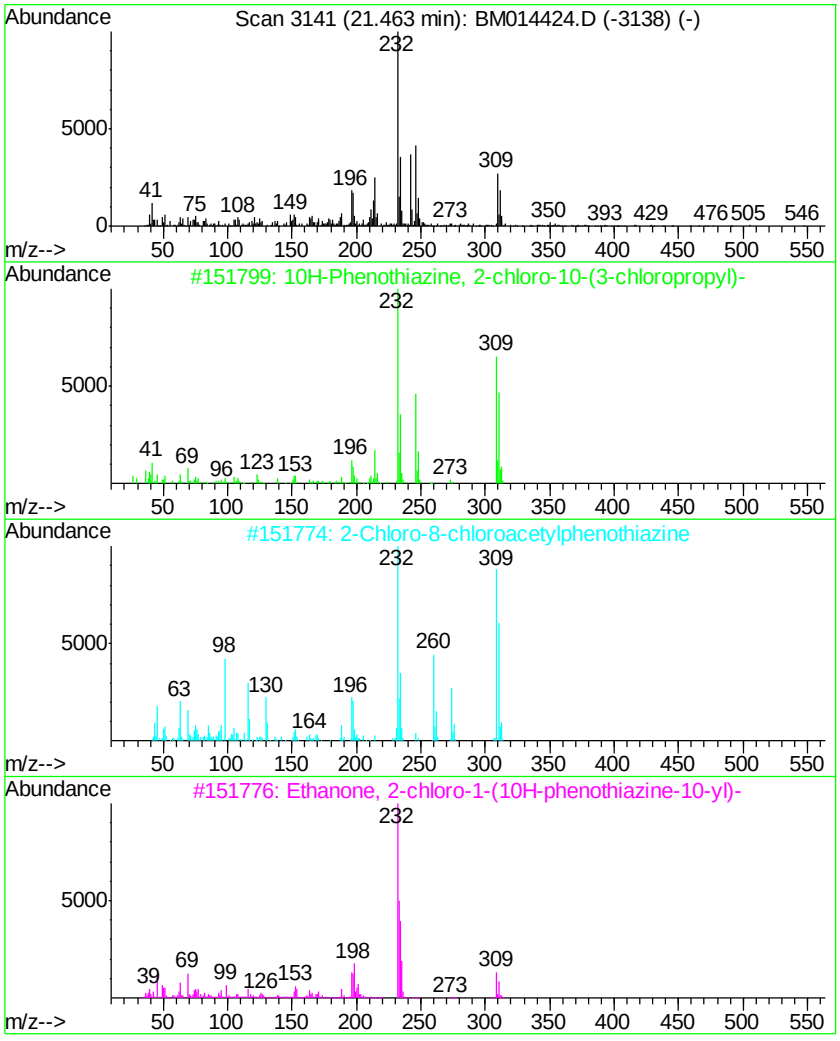
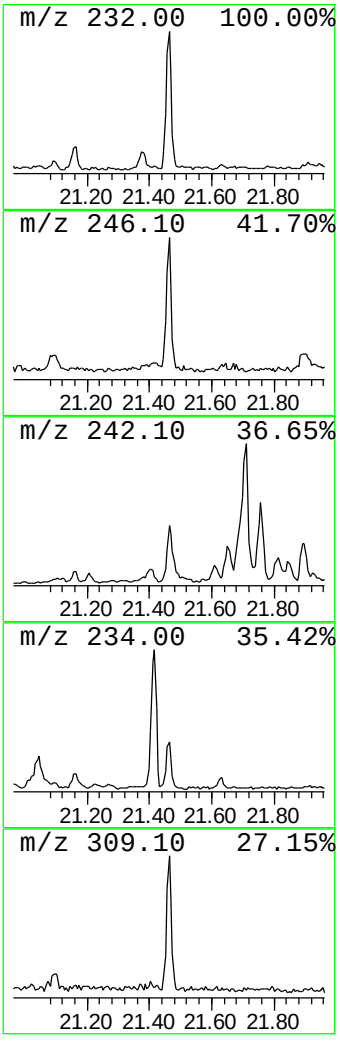
Peak Number 13 10H-Phenothiazine, 2-chloro... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.46	4.38 ng/ul	1451520	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10H-Phenothiazine, 2-chloro-10-(...	309	C15H13Cl2NS	002765-59-5	99
2	2-Chloro-8-chloroacetylphenothia...	309	C14H9Cl2NOS	010554-02-6	60
3	Ethanone, 2-chloro-1-(10H-phenot...	309	C14H9Cl2NOS	016189-69-8	55
4	Pyridine-3-carboxamide, 2-(4-chl...	292	C14H13ClN2O3	309743-02-0	35
5	Phenoxathiin, 10,10-dioxide	232	C12H8O3S	000950-47-0	30





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Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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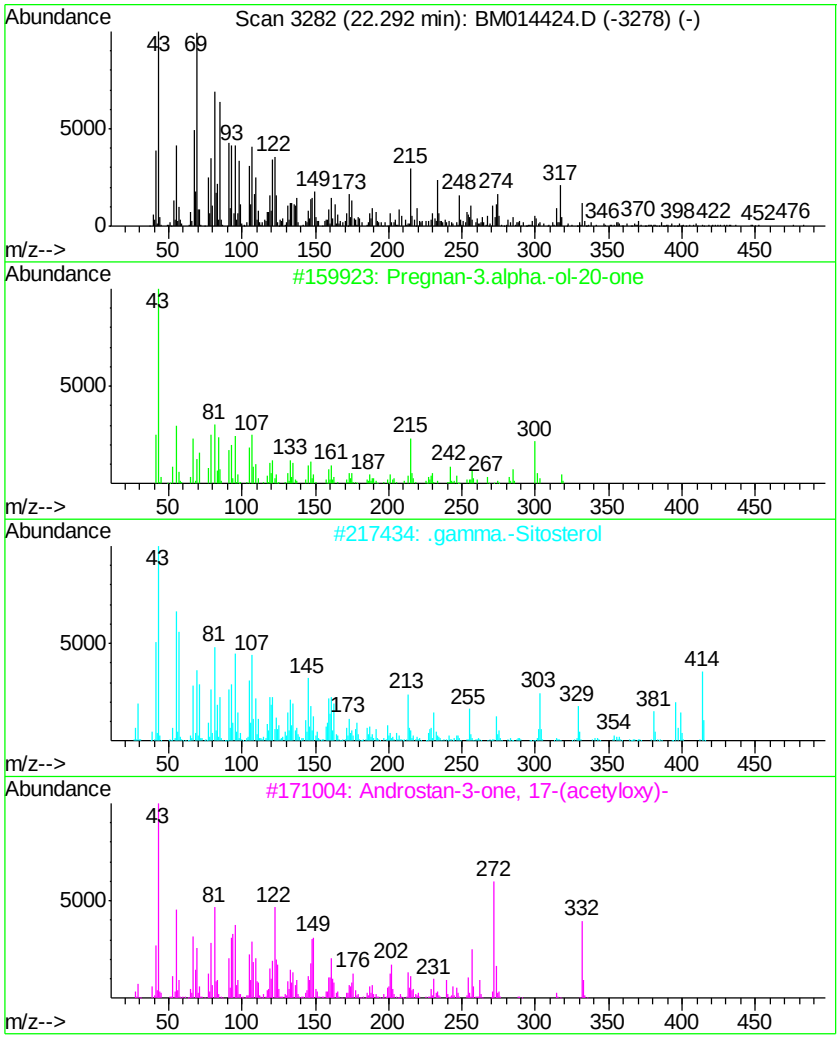
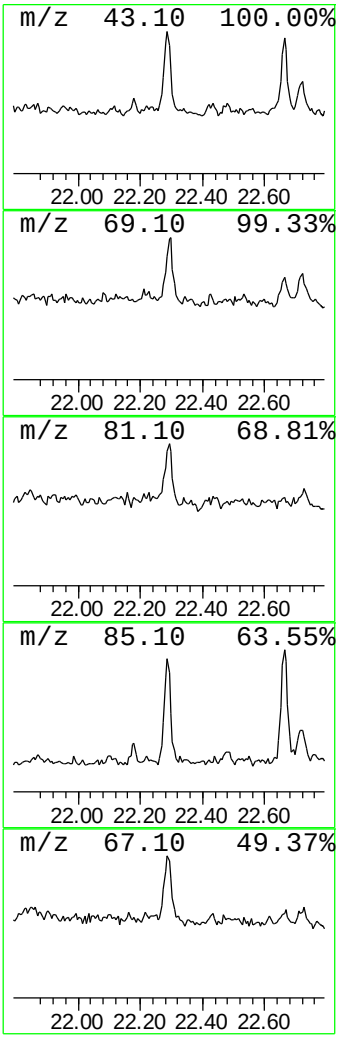
Peak Number 15 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.29	19.50 ng/ul	3073140	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pregnan-3.alpha.-ol-20-one	318	C21H34O2	000128-20-1	46
2	.gamma.-Sitosterol	414	C29H50O	000083-47-6	25
3	Androstan-3-one, 17-(acetyloxy)-	332	C21H32O3	125638-46-2	15
4	Diethylmalonic acid, 1-bromo-3,3...	460	C19H32BrF3O4	1000370-80-1	14
5	4H-1,3-Dioxin-4-one, 2-(1,1-dime...	170	C9H14O3	107289-20-3	11





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Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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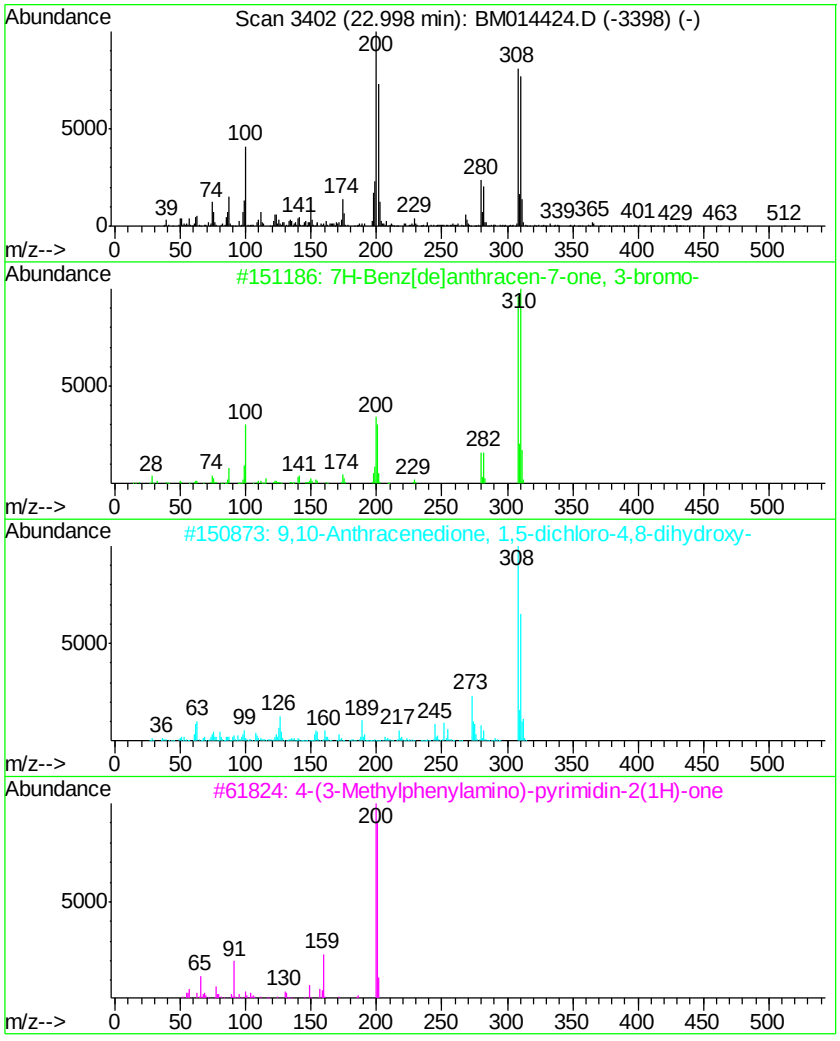
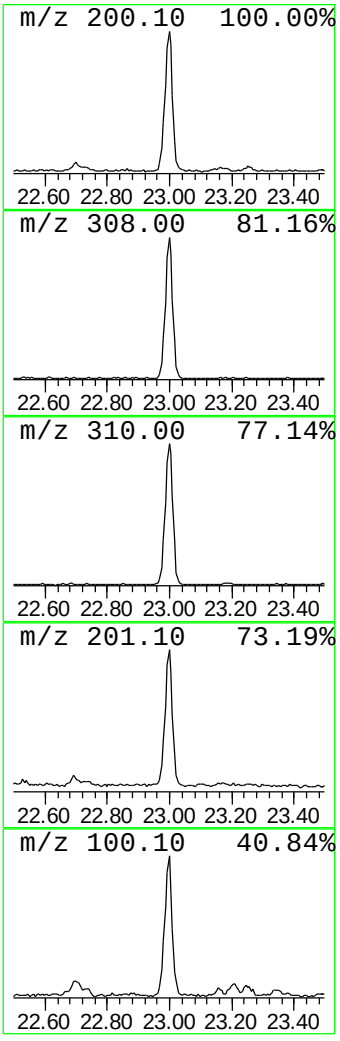
Peak Number 16 7H-Benz[de]anthracen-7-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.00	20.00 ng/ul	3152160	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benz[de]anthracen-7-one, 3-br...	308	C17H9BrO	000081-96-9	94
2	9,10-Anthracenedione, 1,5-dichlo...	308	C14H6Cl2O4	006837-97-4	27
3	4-(3-Methylphenylamino)-pyrimidi...	201	C11H11N3O	127970-30-3	22
4	13-Ethyl-17alpha-ethynyl-17beta-...	308	C21H24O2	1000234-44-4	15
5	4,6-Diamino-1-[3,4-dichloropheny...	308	C12H10Cl2N6	058791-63-2	12





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Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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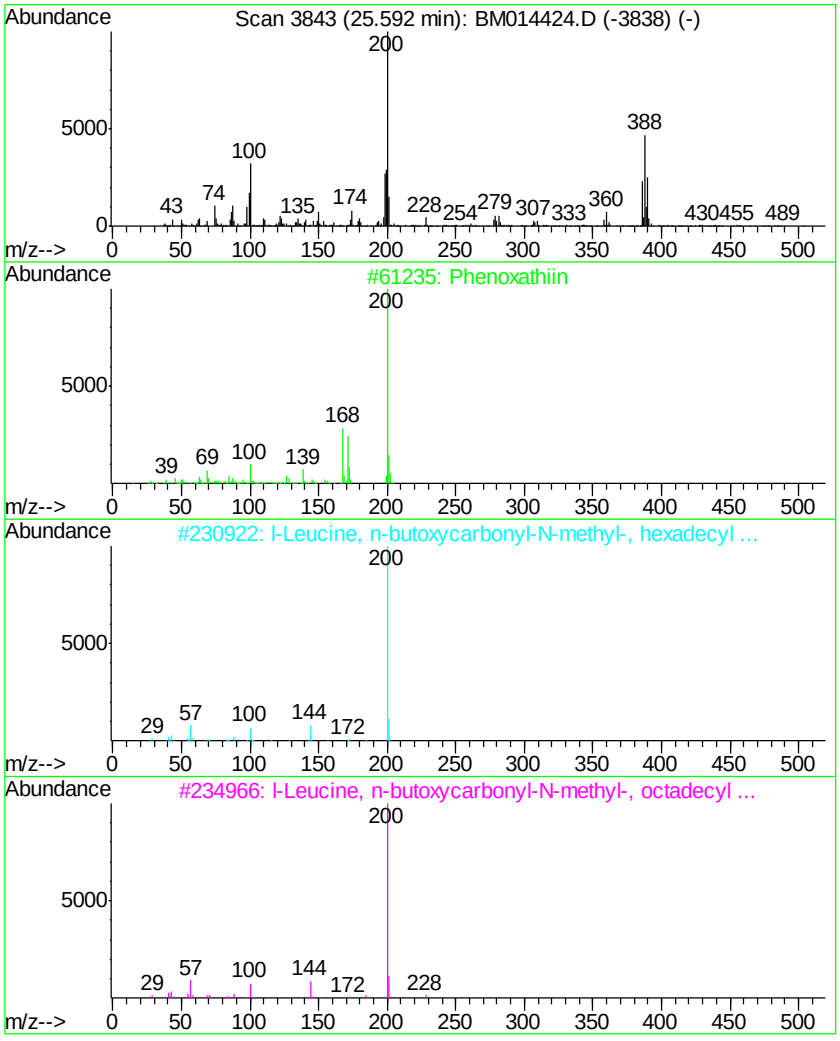
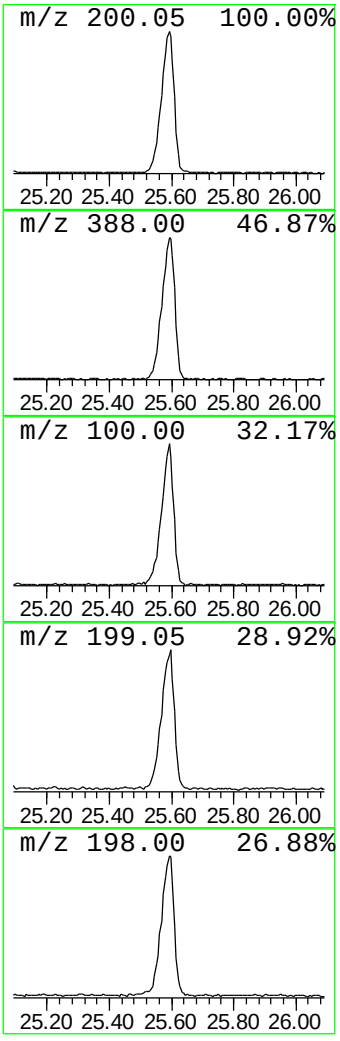
Peak Number 17 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.59	40.16 ng/ul	6329380	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenoxathiin	200	C12H8OS	000262-20-4	25
2	l-Leucine, n-butoxycarbonyl-N-me...	469	C28H55NO4	1000321-89-3	22
3	l-Leucine, n-butoxycarbonyl-N-me...	497	C30H59NO4	1000321-89-5	22
4	D-Norleucine, N-neopentyloxycarb...	455	C27H53NO4	1000344-14-0	14
5	2-Aminocaprylic acid, N-propoxyc...	413	C24H47NO4	1000325-42-9	14





Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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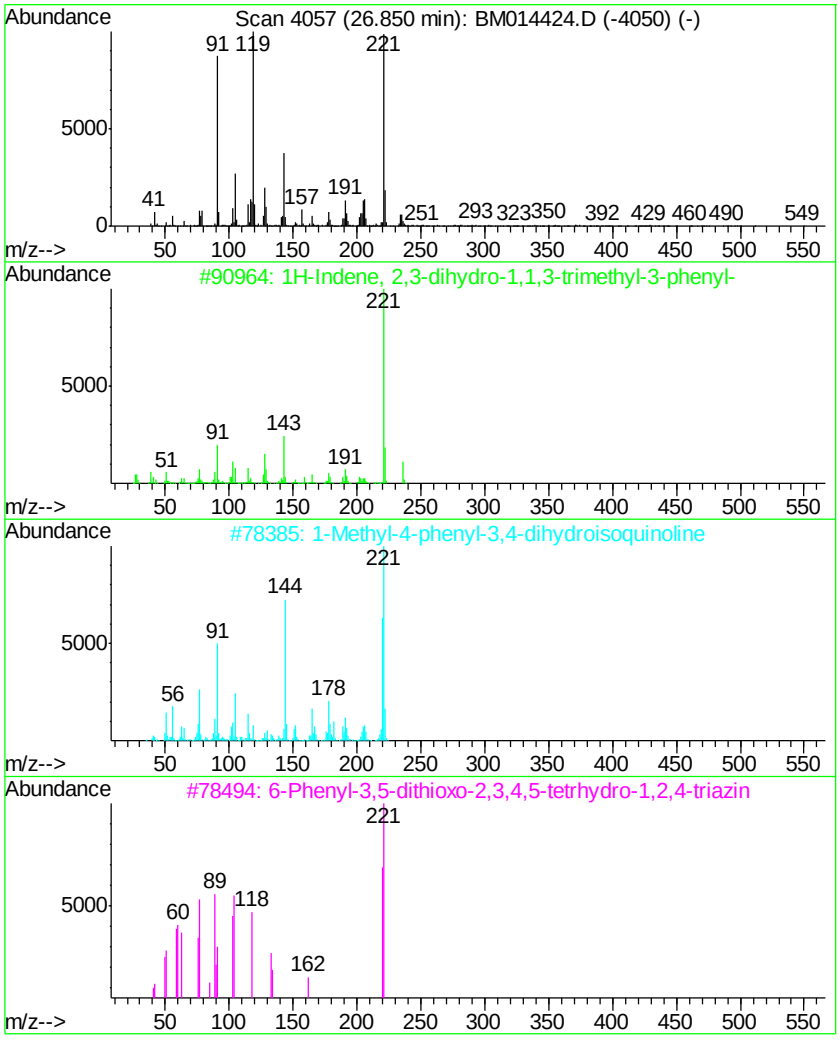
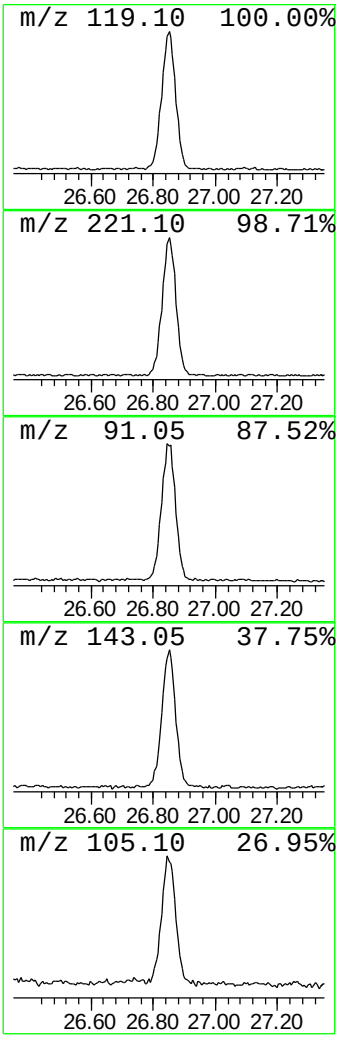
Peak Number 18 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.85	36.35 ng/ul	5728380	Perylene-d12	23.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	43
2	1-Methyl-4-phenyl-3,4-dihydroiso...	221	C16H15N	1000211-01-5	41
3	6-Phenyl-3,5-dithioxo-2,3,4,5-te...	221	C9H7N3S2	038119-57-2	35
4	Benz[j]isoquinoline, 3,6,8-trime...	221	C16H15N	061171-16-2	30
5	(.+/-.)-2-Phenylbutyric acid, te...	278	C16H26O2Si	1000333-14-6	27





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022818\
Data File : BM014424.D
Acq On : 01 Mar 2018 05:09
Operator : SJ/JU
Sample : J1646-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE5W8

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Phenanthrene, 2-m...	17.86	14.0	ng/ul	1658620	4	16.93	2375210	20.0
9,10-Anthracenedione	18.37	16.2	ng/ul	1919400	4	16.93	2375210	20.0
1,1'-Biphenyl, 2,...	19.15	4.5	ng/ul	1481170	5	21.16	6630290	20.0
Mitotane	20.01	36.1	ng/ul	11963200	5	21.16	6630290	20.0
10H-Phenothiazine...	20.20	7.4	ng/ul	2465000	5	21.16	6630290	20.0
3-(4-Methoxybenzy...	20.23	12.7	ng/ul	4221430	5	21.16	6630290	20.0
(DEL) Alkane: Cyc...	20.87	6.3	ng/ul	2073080	5	21.16	6630290	20.0
7H-Benz[de]anthra...	21.37	16.4	ng/ul	5423700	5	21.16	6630290	20.0
1H-Indene, 2,3-di...	21.42	76.3	ng/ul	25306700	5	21.16	6630290	20.0
10H-Phenothiazine...	21.46	4.4	ng/ul	1451520	5	21.16	6630290	20.0
unknown-01	22.29	19.5	ng/ul	3073140	6	23.34	3152160	20.0
7H-Benz[de]anthra...	23.00	20.0	ng/ul	3152160	6	23.34	3152160	20.0
unknown-02	25.59	40.2	ng/ul	6329380	6	23.34	3152160	20.0
unknown-03	26.85	36.4	ng/ul	5728380	6	23.34	3152160	20.0