LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100418\
Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.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	6.816	644	651	665	rVB	403727	770472	1.56%	0.535%
2	7.163	704	710	720	rBV	425841	692764	1.40%	0.481%
3	7.622	781	788	798	rBV	498204	805119	1.63%	0.559%
4	8.340	902	910	920	rBV2	435198	785441	1.59%	0.545%
5	8.775	977	984	993	rBV	396360	674298	1.37%	0.468%
6	9.487	1099	1105	1115	rBV	336029	583000	1.18%	0.405%
7	10.034	1191	1198	1213	rBV	491599	939113	1.90%	0.652%
8	10.393	1252	1259	1264	rBV	743153	1245202	2.52%	0.864%
9	10.540	1278	1284	1306	rBV	332796	674992	1.37%	0.468%
10	13.675	1812	1817	1822	rBV	995000	1360953	2.76%	0.945%
11	13.951	1855	1864	1877	rBV	1214210	1826188	3.70%	1.267%
12	14.263	1910	1917	1923	rBV2	1101865	1684881	3.42%	1.169%
13	14.504	1949	1958	1968	rBV	275689	592804	1.20%	0.411%
14	15.263	2080	2087	2092	rBV	1508412	2236607	4.53%	1.552%
15	15.404	2106	2111	2121	rBV	396038	654858	1.33%	0.454%
16	17.016	2377	2385	2388	rBV	1389718	2055653	4.17%	1.427%
17	17.057	2388	2392	2397	rVV	2910082	4032805	8.18%	2.799%
18	17.116	2397	2402	2405	rVV	1767161	2529595	5.13%	1.756%
19	17.145	2405	2407	2413	rVB	542169	702393	1.42%	0.487%
20	17.422	2445	2454	2461	rBV2	425107	716075	1.45%	0.497%
21	17.939	2536	2542	2547	rVB2	419452	864843	1.75%	0.600%
22	18.086	2561	2567	2571	rBV2	480634	788081	1.60%	0.547%
23	19.086	2730	2737	2744	rVB	5364669	7184560	14.57%	4.986%
24	19.422	2788	2794	2796	rBV2	2616054	4119047	8.35%	2.859%
25	19.451	2796	2799	2807	rVB	4677295	5994230	12.15%	4.160%
26	19.780	2848	2855	2861	rBV	6779430	8351968	16.93%	5.797%
27	19.945	2879	2883	2889	rVB2	611328	885129	1.79%	0.614%
28	20.045	2896	2900	2904	rBV	560800	719223	1.46%	0.499%
29	20.104	2904	2910	2914	rBV2	295346	495511	1.00%	0.344%
30	20.898	3042	3045	3048	rVV	488512	627194	1.27%	0.435%
31	20.939	3048	3052	3058	rVV3	664319	1327163	2.69%	0.921%
32	21.157	3082	3089	3093	rBV	32017377	49323293	100.00%	34.232%
33	21.210	3093	3098	3103	rVV2	2825879	5965892	12.10%	4.140%
34	21.257	3103	3106	3116	rVB	2562078	3570560	7.24%	2.478%

LSC Area Percent Report

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Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Title : SVOA CALIBRATION

35	21.374	3123	3126	3131	rVB2	685981	852504	1.73%	0.592%
36	21.721	3181	3185	3188	rBV3	597958	664962	1.35%	0.462%
37	21.810	3196	3200	3206	rBV3	342854	582395	1.18%	0.404%
38	21.874	3206	3211	3216	rVB4	1279320	1932354	3.92%	1.341%
39	21.963	3223	3226	3230	rBV3	330612	572327	1.16%	0.397%
40	22.027	3233	3237	3239	rVV3	1040306	1602249	3.25%	1.112%
41	22.051	3239	3241	3249	rVB2	1123242	1508562	3.06%	1.047%
42	22.210	3265	3268	3274	rVB4	876748	1558316	3.16%	1.082%
43	22.386	3293	3298	3301	rBV2	475588	848809	1.72%	0.589%
44	22.774	3356	3364	3369	rBV2	1874094	4873721	9.88%	3.382%
45	23.239	3438	3443	3447	rBV	947608	1721246	3.49%	1.195%
46	23.292	3447	3452	3456	rVV	1173443	2376960	4.82%	1.650%
47	23.339	3456	3460	3467	rVB	1541091	2562936	5.20%	1.779%
48	23.427	3470	3475	3479	rVV	848359	1555752	3.15%	1.080%
49	23.474	3480	3483	3491	rVB	462162	797664	1.62%	0.554%
50	23.927	3556	3560	3572	rVB	278245	615411	1.25%	0.427%
51	25.627	3841	3849	3859	rVB	732508	2060199	4.18%	1.430%
52	26.298	3955	3963	3972	rBV	573884	1620092	3.28%	1.124%

Sum of corrected areas: 144086366

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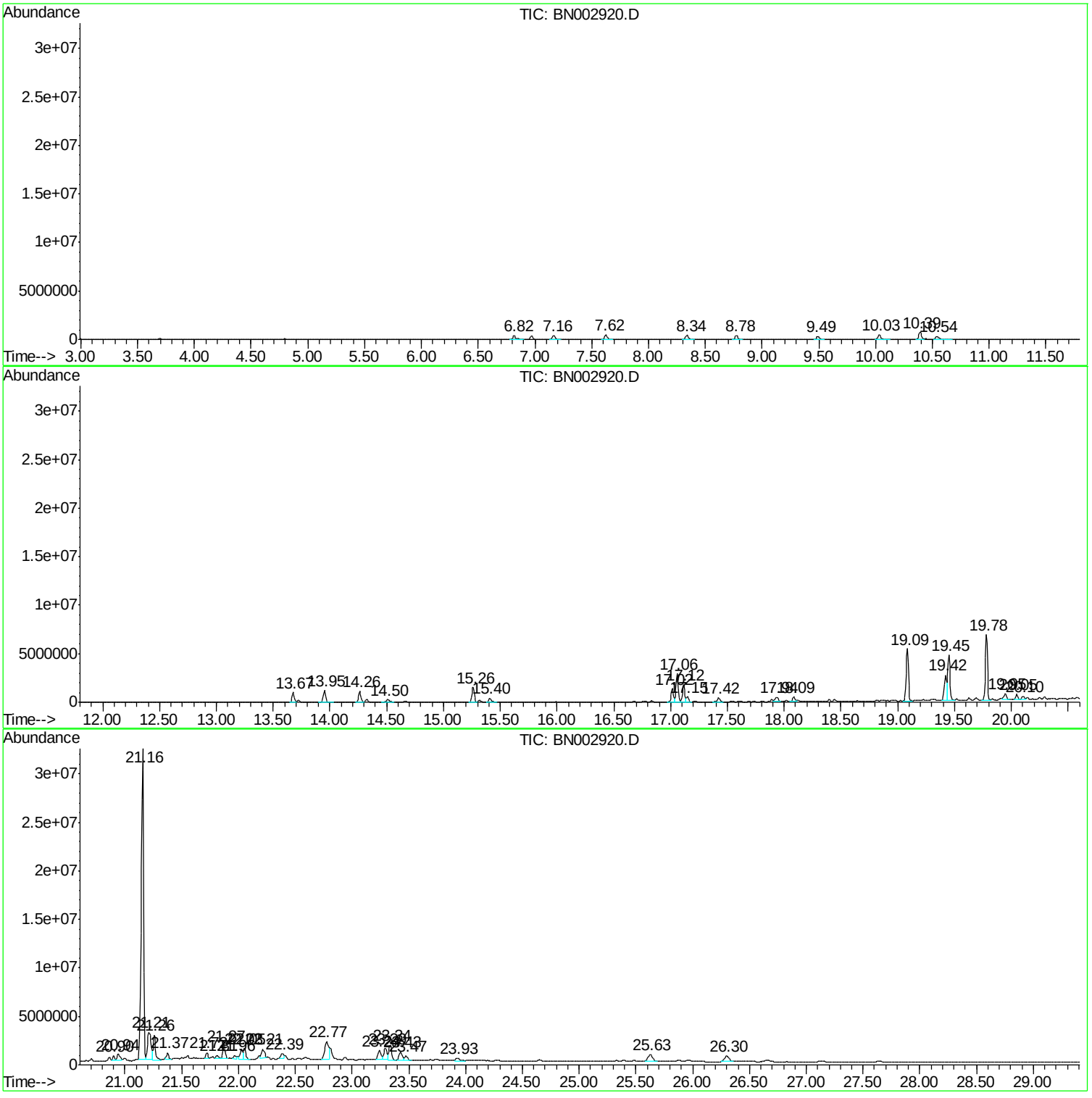
Instrument :
BNA_N
ClientSampleId :
JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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



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Quant Title : SVOA CALIBRATION

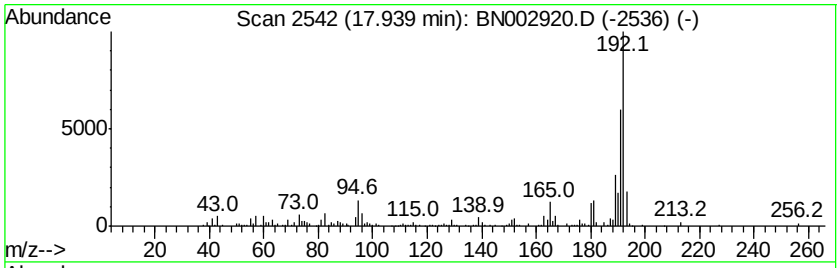
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

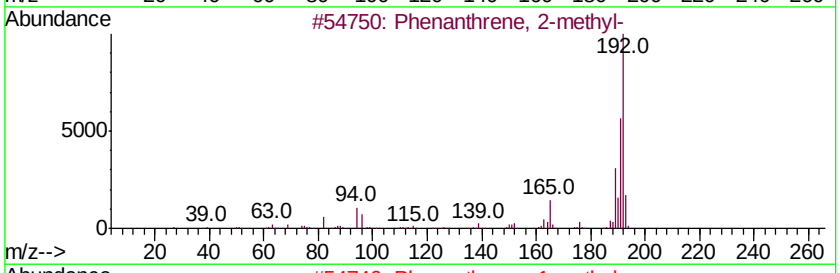
****

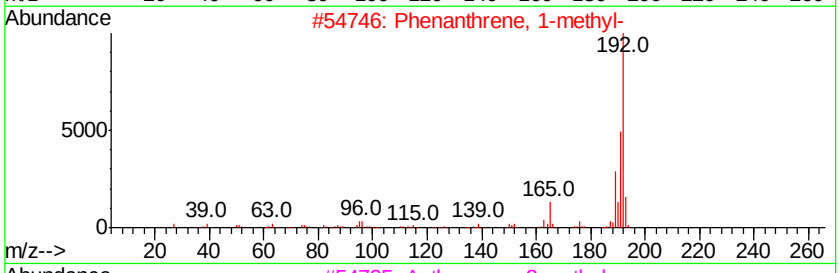
Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 4

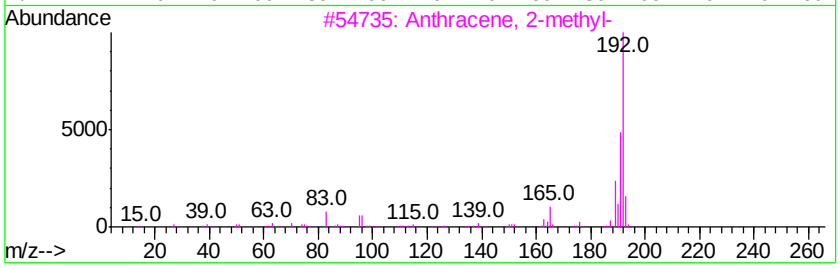
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	8.41 ng/ul	864843	Phenanthrene-d10	17.02

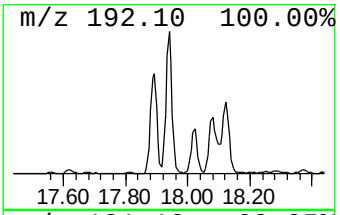
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94

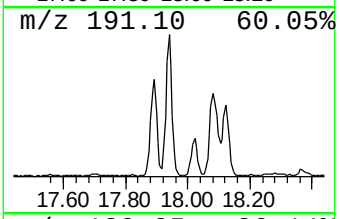


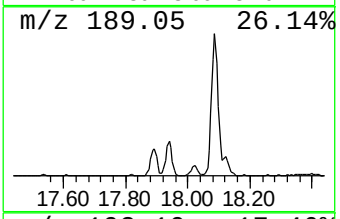


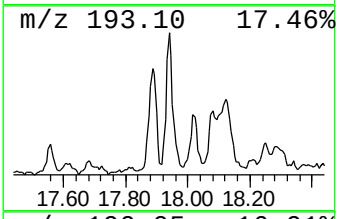


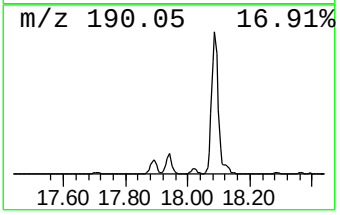












Library Search Compound Report

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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

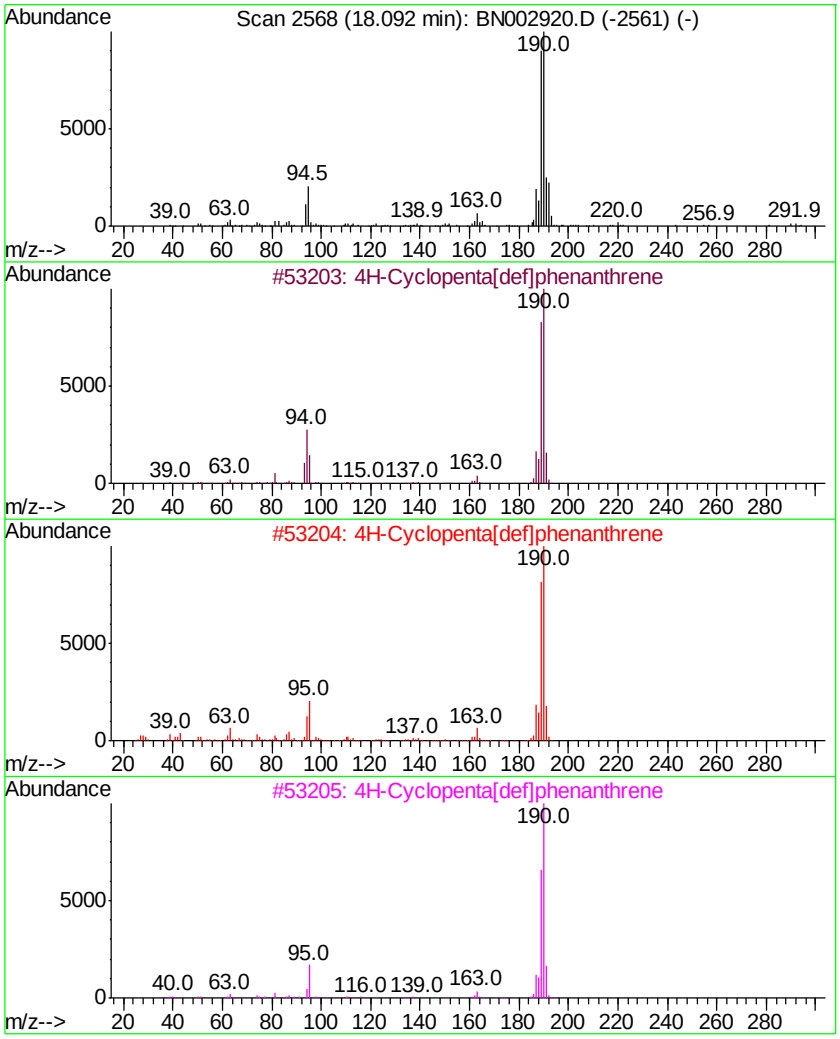
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

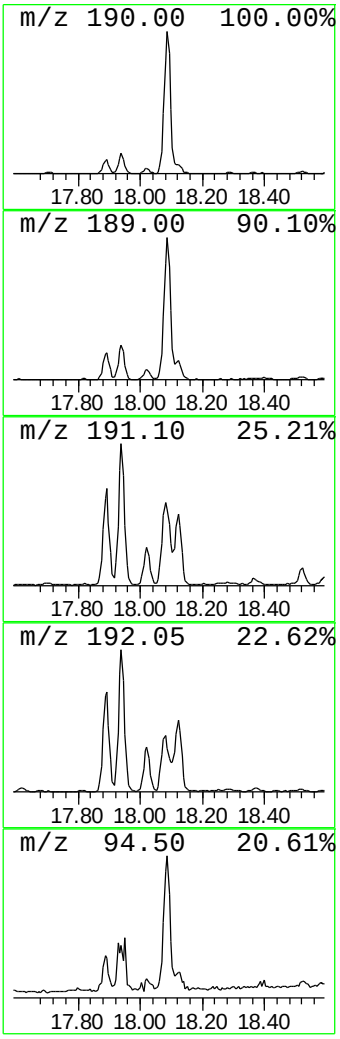
****

Peak Number 2 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	7.67 ng/ul	788081	Phenanthrene-d10	17.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	38





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Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

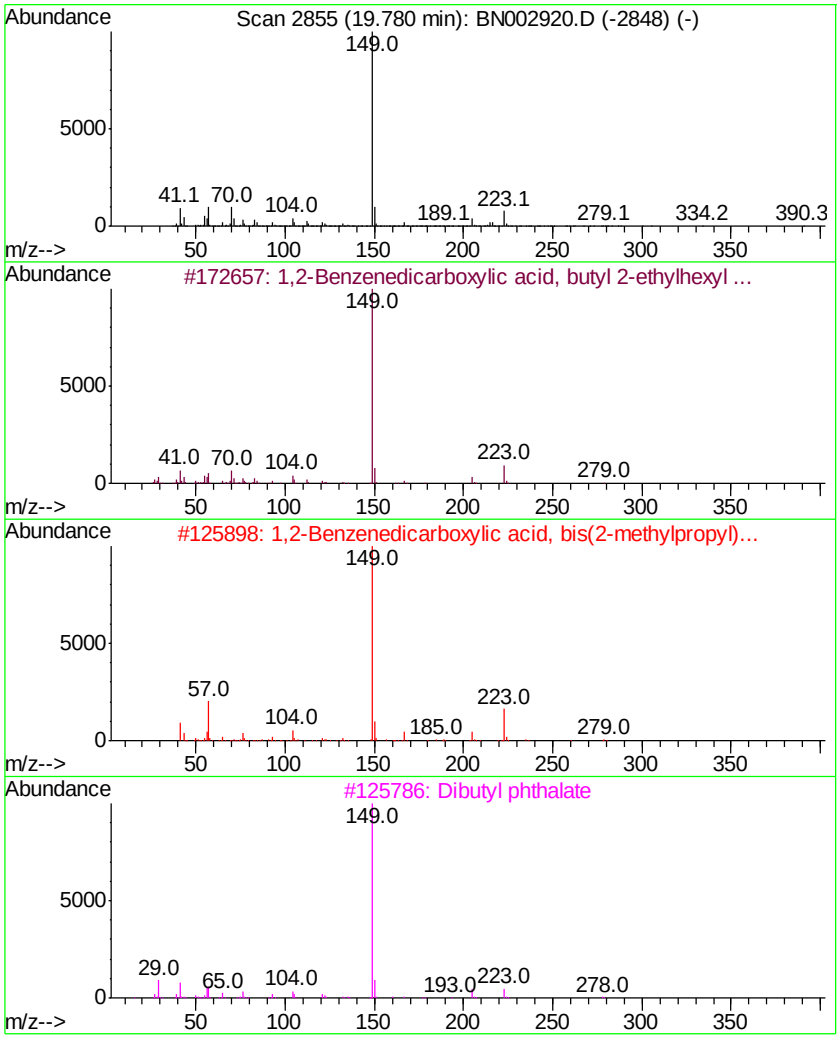
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

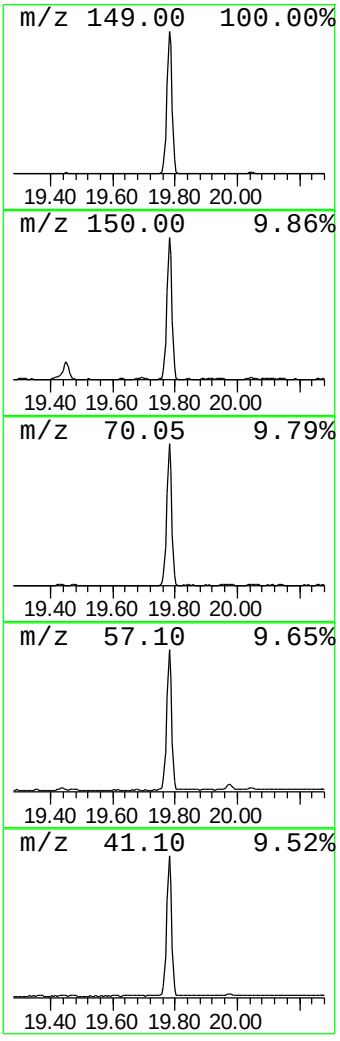
****

Peak Number 3 1,2-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.78	28.00 ng/ul	8351970	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	91
2	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	90
3	Dibutyl phthalate	278	C16H22O4	000084-74-2	86
4	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000084-78-6	86
5	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000084-78-6	80





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Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

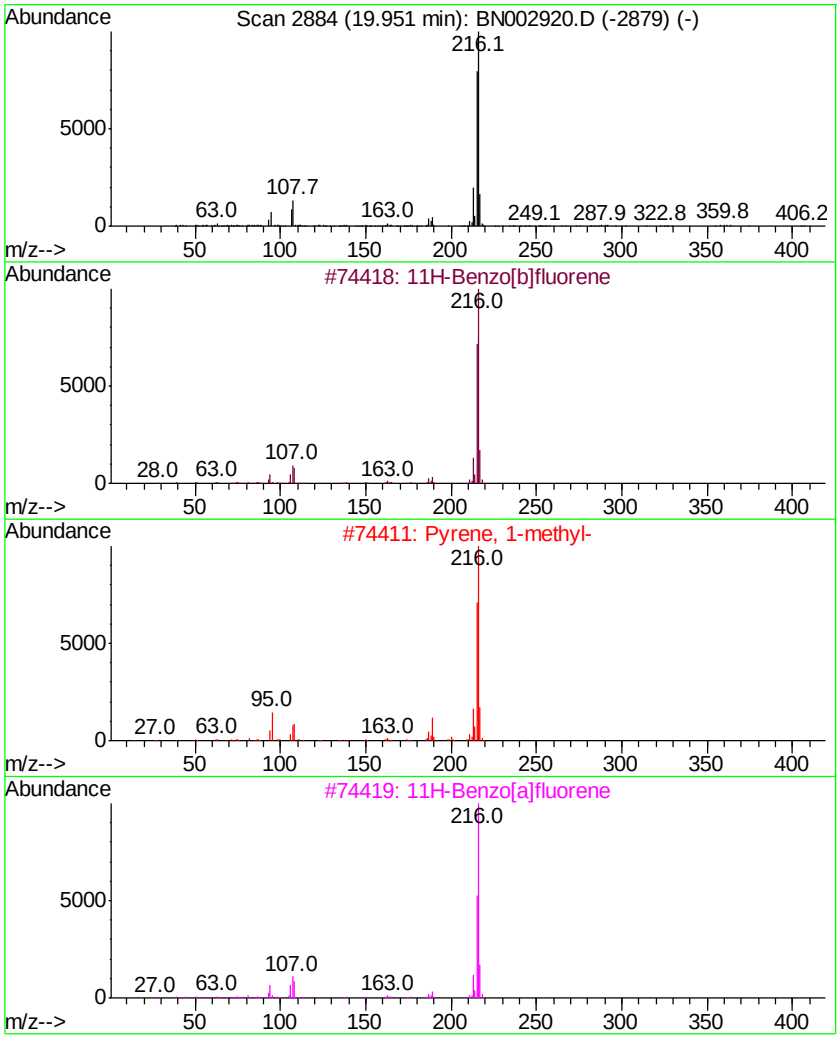
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

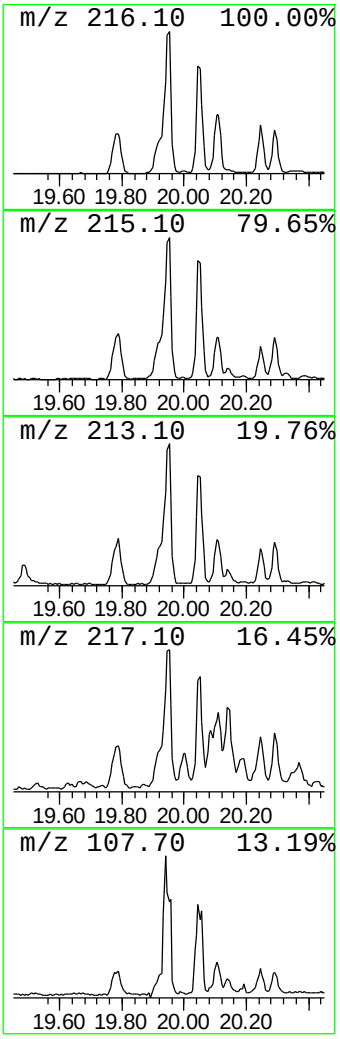
****

Peak Number 4 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.95	2.97 ng/ul	885129	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91





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Instrument : BNA_N
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

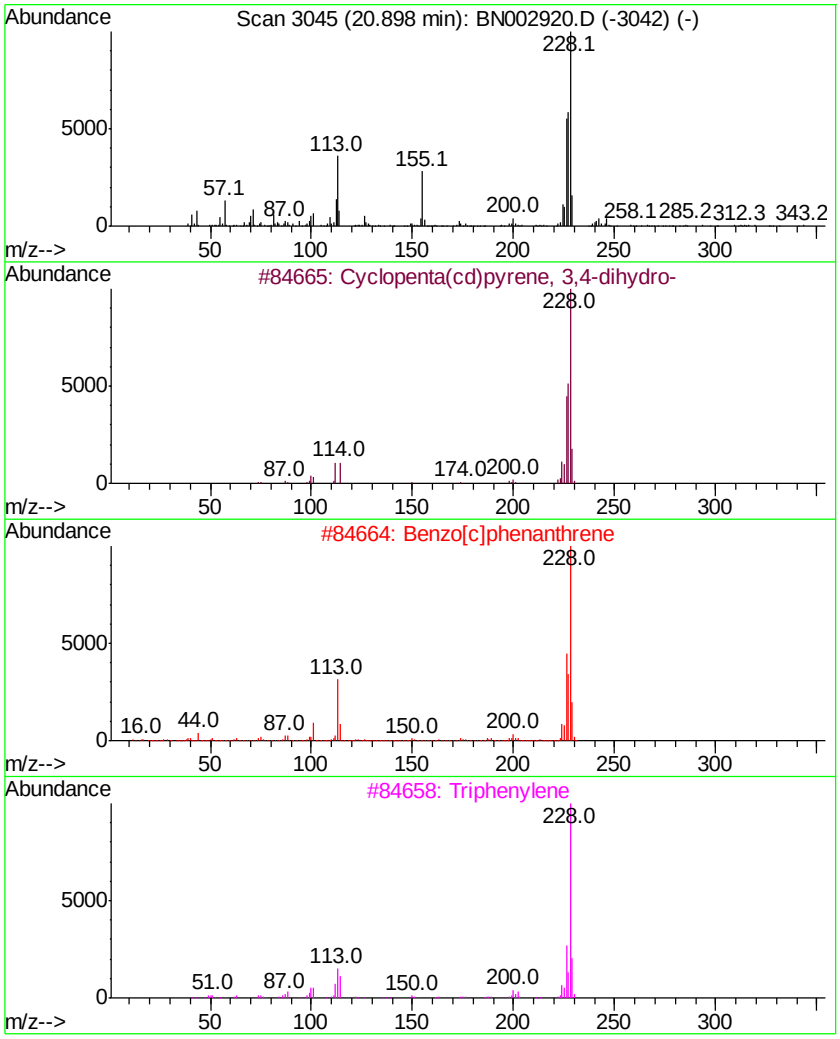
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

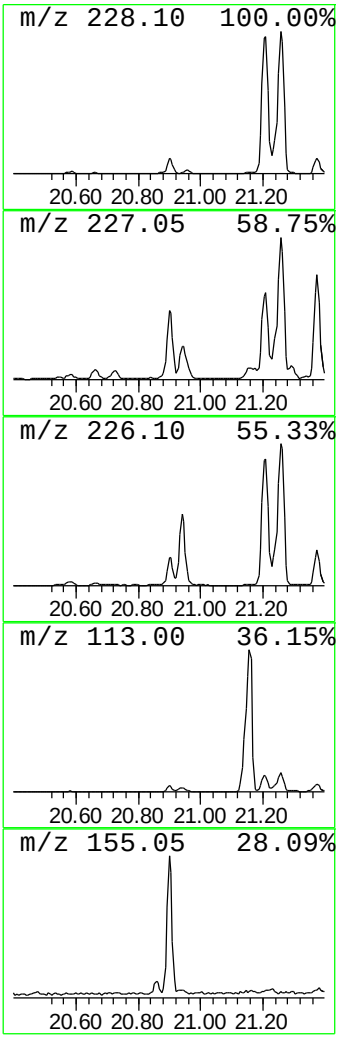
****

Peak Number 6 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.90	2.10 ng/ul	627194	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	70
2	Benzo[c]phenanthrene	228	C18H12	000195-19-7	60
3	Triphenylene	228	C18H12	000217-59-4	50
4	Benzo[c]phenanthrene	228	C18H12	000195-19-7	47
5	Benz[a]anthracene	228	C18H12	000056-55-3	41





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

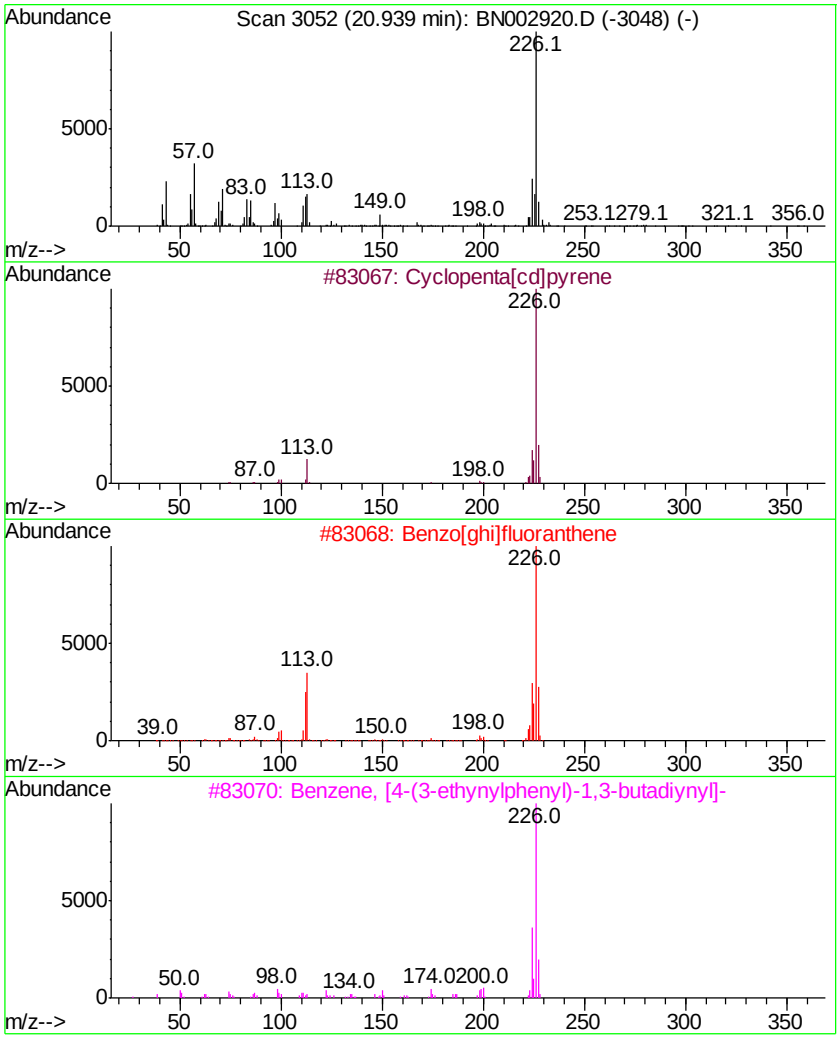
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

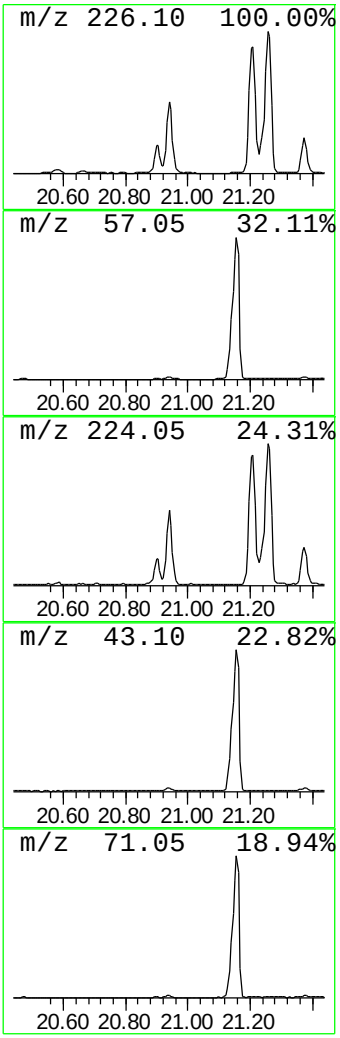
****

Peak Number 7 Cyclopenta[cd]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	4.45 ng/ul	1327160	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	60
2	Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	58
3	Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	38
4	6-Bromo-.alpha.-[di-n-heptylamin...	511	C30H42BrNO	027022-39-5	32
5	9-Methoxymethyl-1-methyl-3,6-dia...	226	C12H22N2O2	1000216-30-2	32





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Instrument : BNA_N
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Quant Title : SVOA CALIBRATION

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

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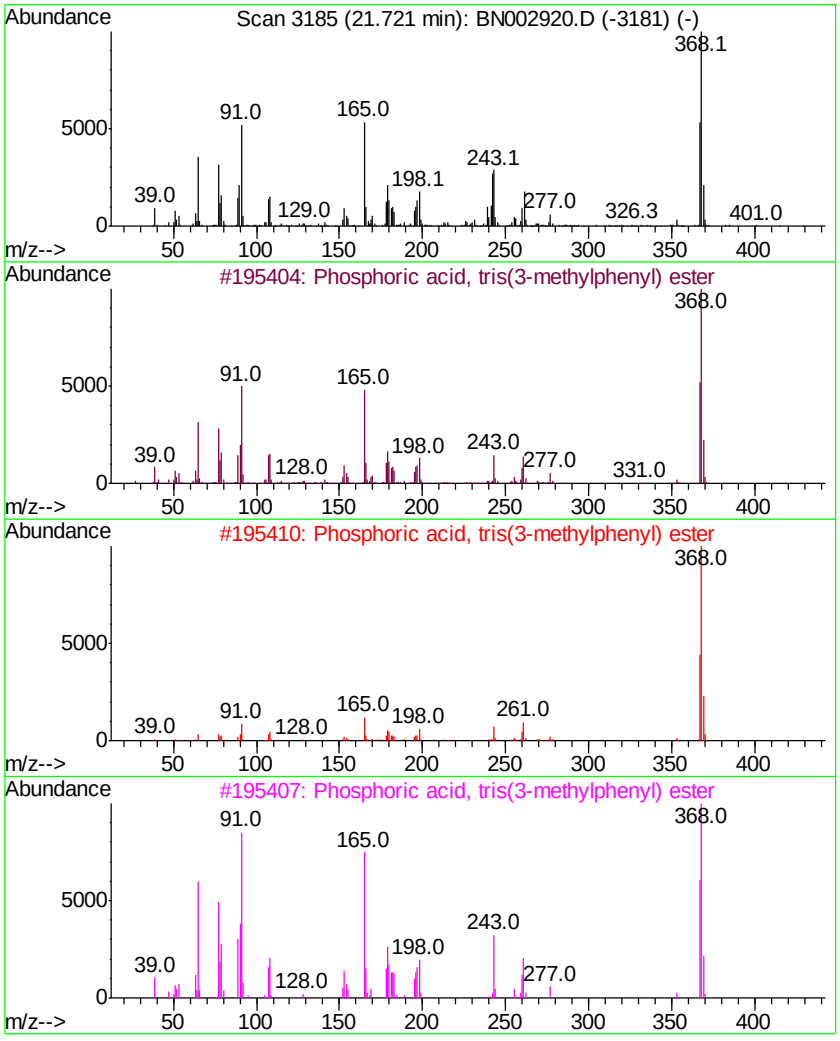
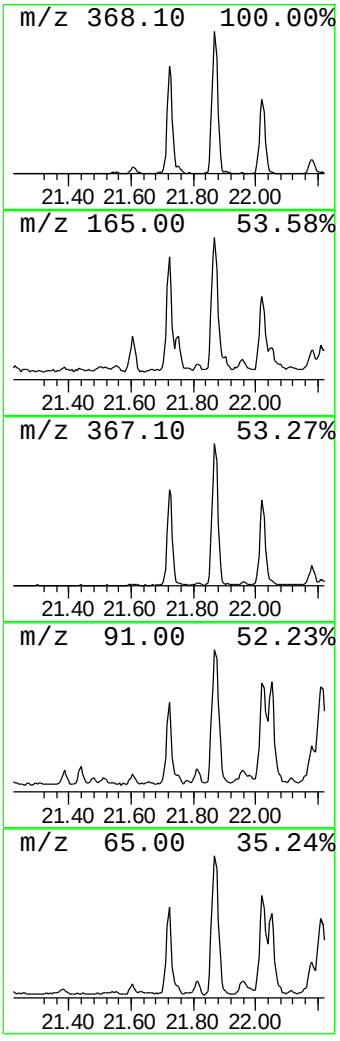
Peak Number 9 Phosphoric acid, tris(3-met... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	2.23 ng/ul	664962	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	99
2	Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	99
3	Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	99
4	Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	94
5	Phosphoric acid, tris(2-methylph...	368	C21H21O4P	000078-30-8	81





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100418\
Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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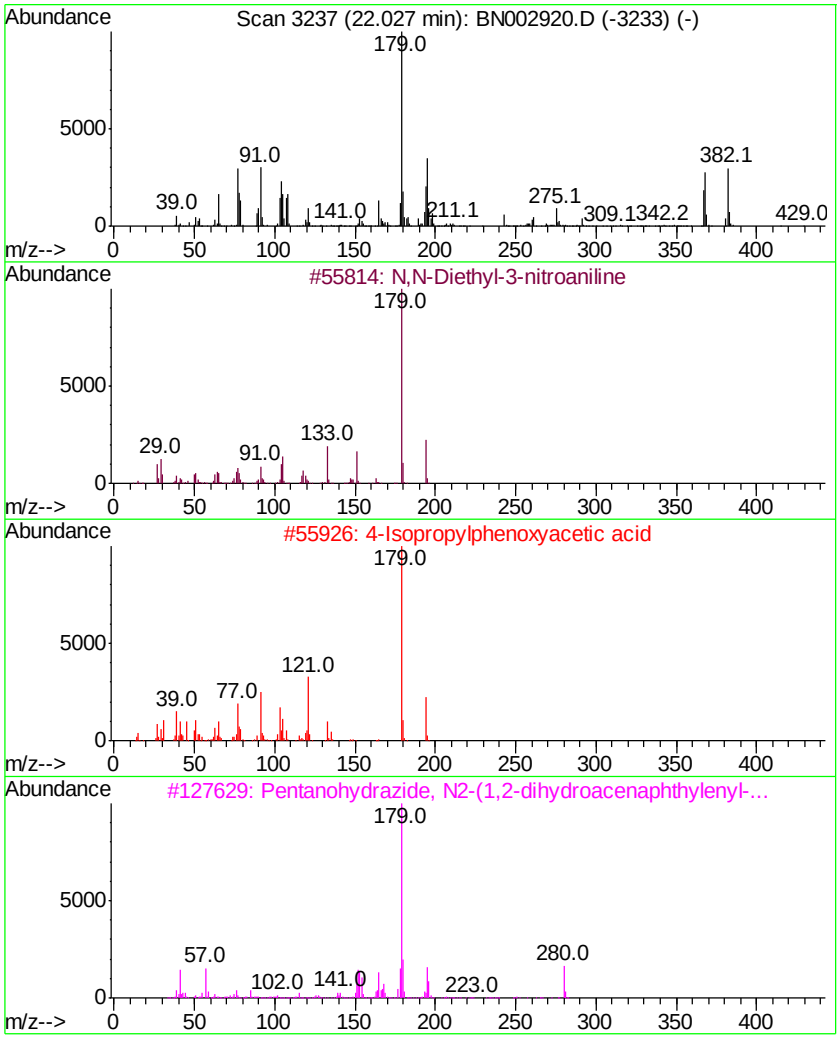
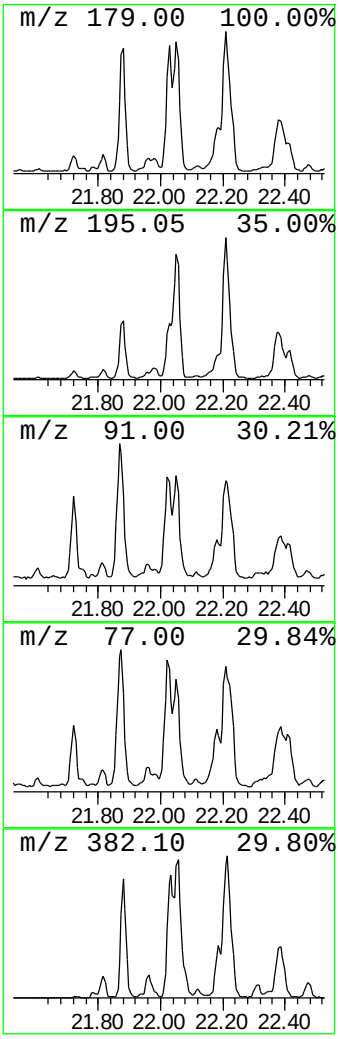
Peak Number 11 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.03	5.37 ng/ul	1602250	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N,N-Diethyl-3-nitroaniline	194	C10H14N2O2	002216-16-2	38
2	4-Isopropylphenoxyacetic acid	194	C11H14O3	001643-16-9	38
3	Pentanohydrazide, N2-(1,2-dihydr...	280	C18H20N2O	341938-83-8	25
4	Nicotinamide N-oxide, N-tert-but...	252	C12H20N2O2Si	1000364-74-4	23
5	1H-1,2,3,4-Tetrazole-1-acetamide...	221	C9H8FN5O	1000338-06-1	22





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Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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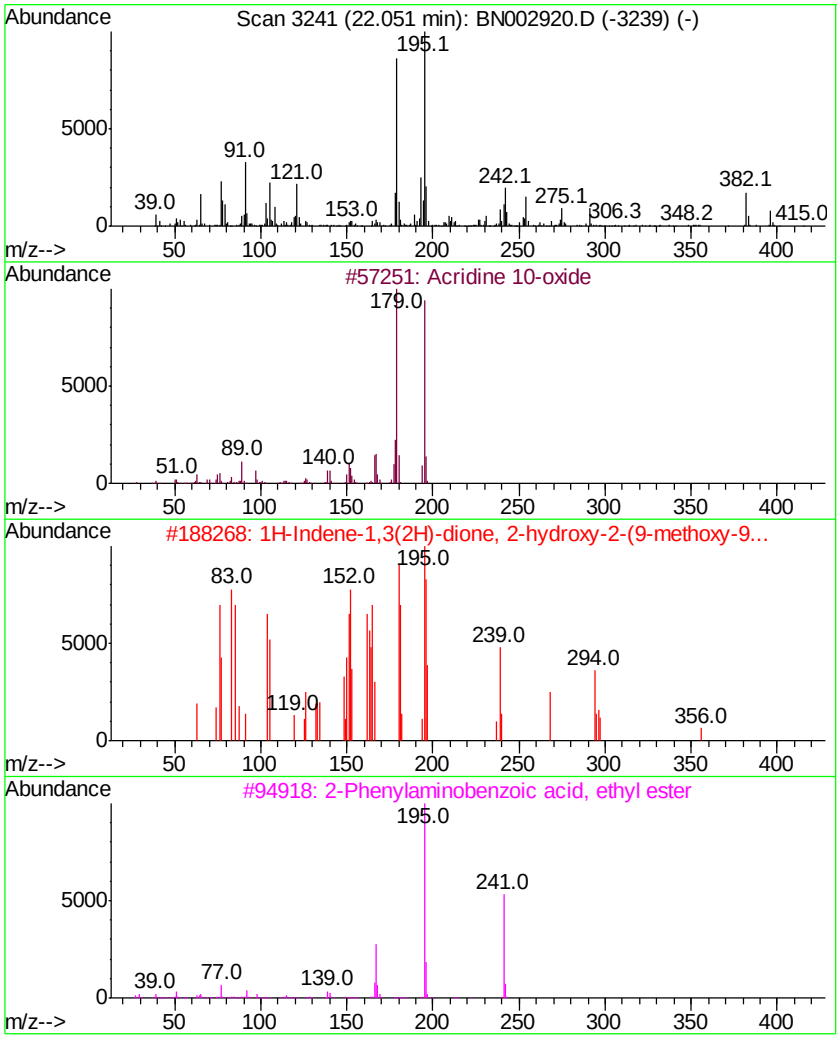
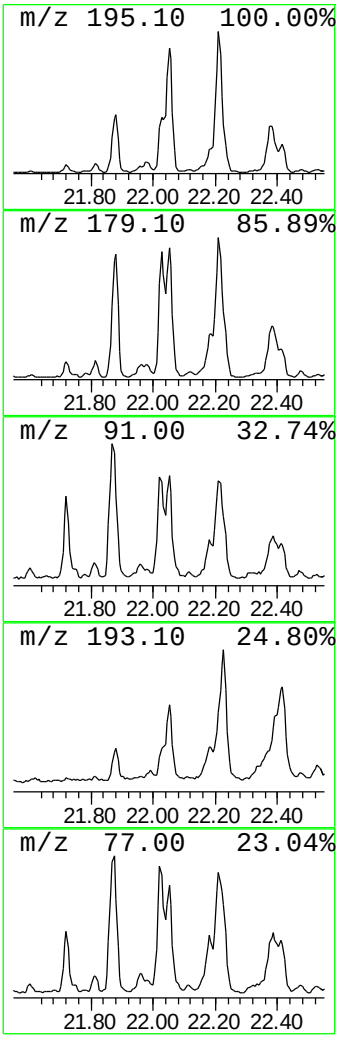
Peak Number 12 Acridine 10-oxide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.05	5.06 ng/ul	1508560	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridine 10-oxide	195	C13H9NO	010399-73-2	58
2	1H-Indene-1,3(2H)-dione, 2-hydro...	356	C23H16O4	050616-96-1	35
3	2-Phenylaminobenzoic acid, ethyl...	241	C15H15NO2	023868-11-3	27
4	Carbazole, 1,6-dimethyl-	195	C14H13N	078787-77-6	27
5	2-Phenazinamine	195	C12H9N3	002876-23-5	27





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Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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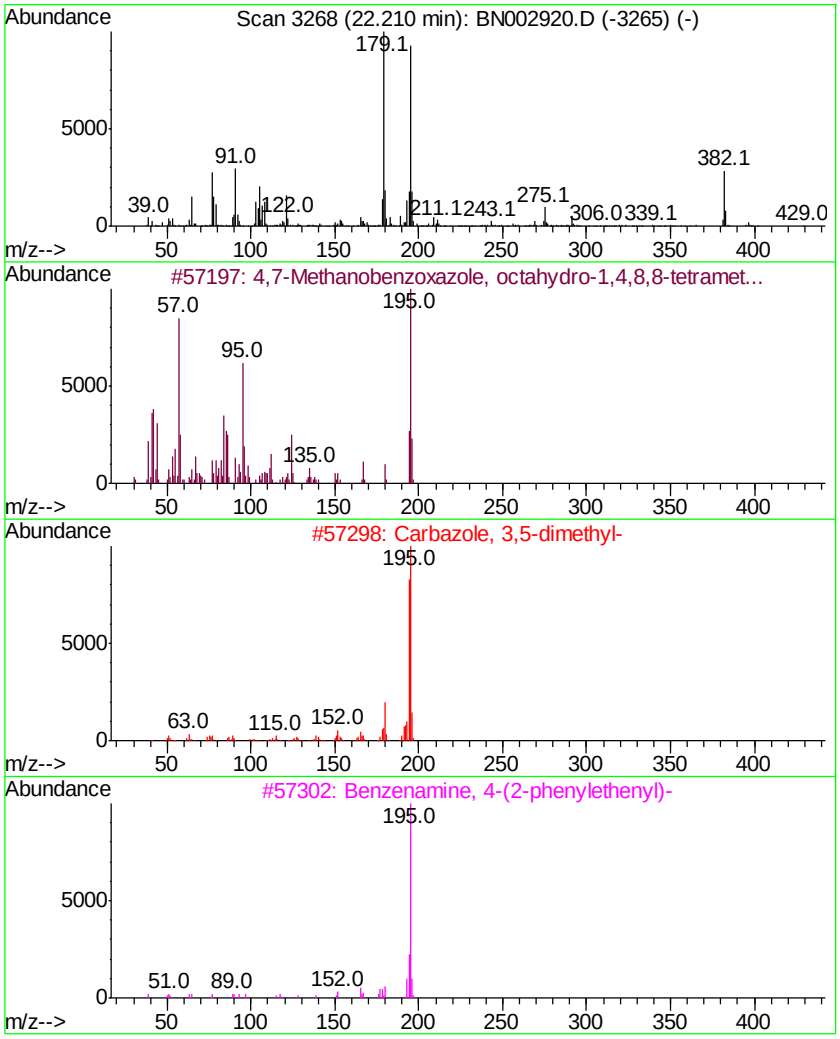
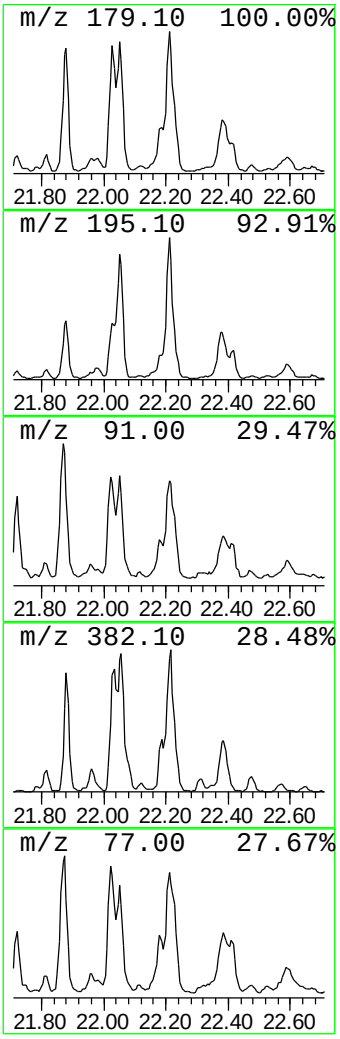
Peak Number 13 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.21	5.22 ng/ul	1558320	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,7-Methanobenzoxazole, octahydr...	195	C12H21NO	035972-97-5	42
2	Carbazole, 3,5-dimethyl-	195	C14H13N	078787-81-2	38
3	Benzenamine, 4-(2-phenylethenyl)-	195	C14H13N	000834-24-2	38
4	Carbazole, 1,5-dimethyl-	195	C14H13N	051640-60-9	38
5	Carbazole, 4,5-dimethyl-	195	C14H13N	018992-66-0	38





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100418\
Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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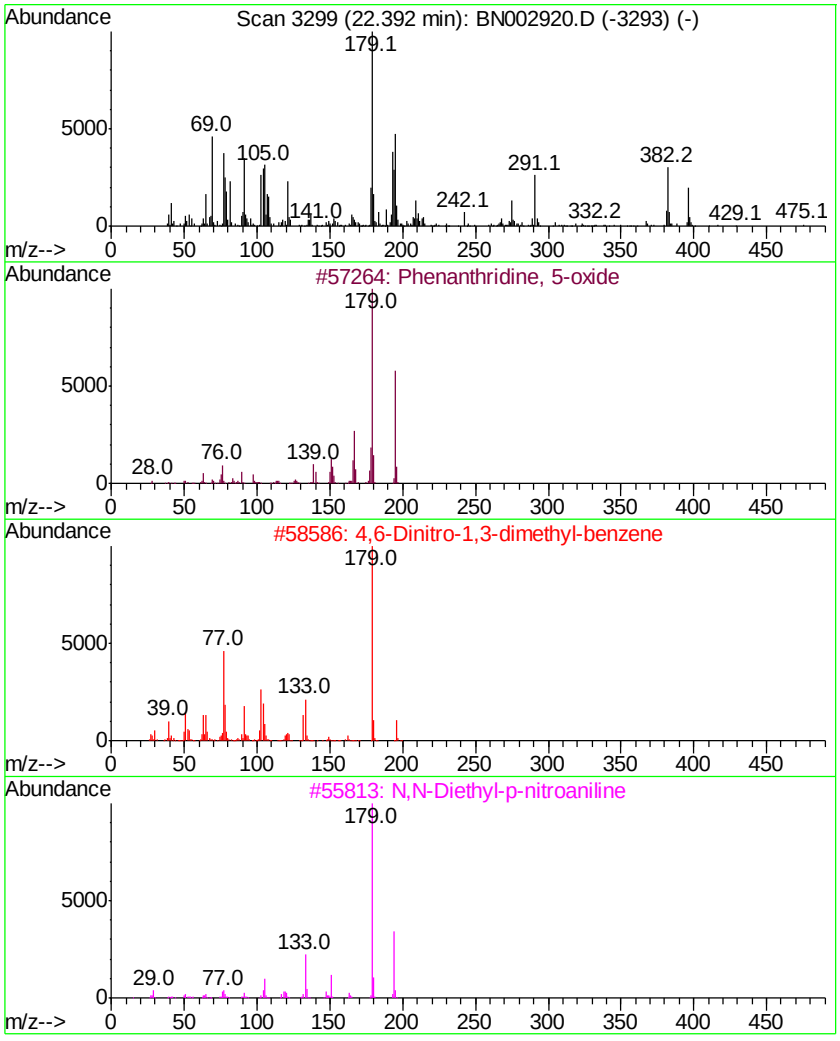
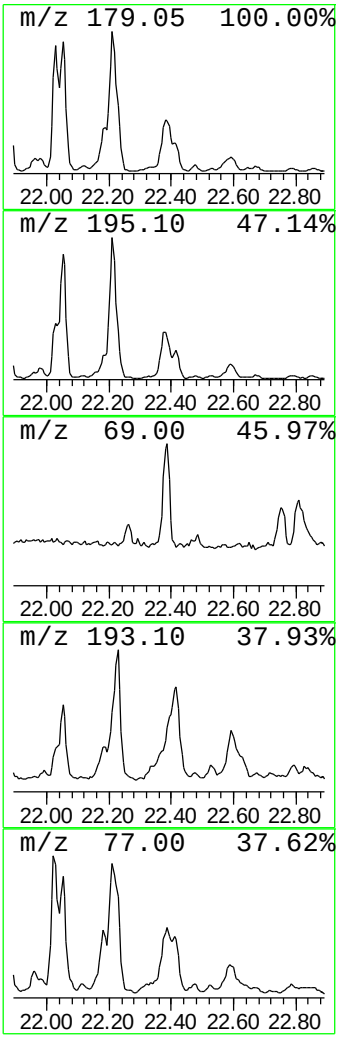
Peak Number 14 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.39	10.91 ng/ul	848809	Perylene-d12	23.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthridine, 5-oxide	195	C13H9NO	014548-01-7	38
2	4,6-Dinitro-1,3-dimethyl-benzene	196	C8H8N2O4	000616-72-8	30
3	N,N-Diethyl-p-nitroaniline	194	C10H14N2O2	002216-15-1	30
4	4-tert-Butylcatechol, dimethyl e...	194	C12H18O2	1000352-79-6	27
5	(2,4-Dimethoxyphenyl)carbamic ac...	291	C15H14FNO4	1000305-11-7	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100418\
Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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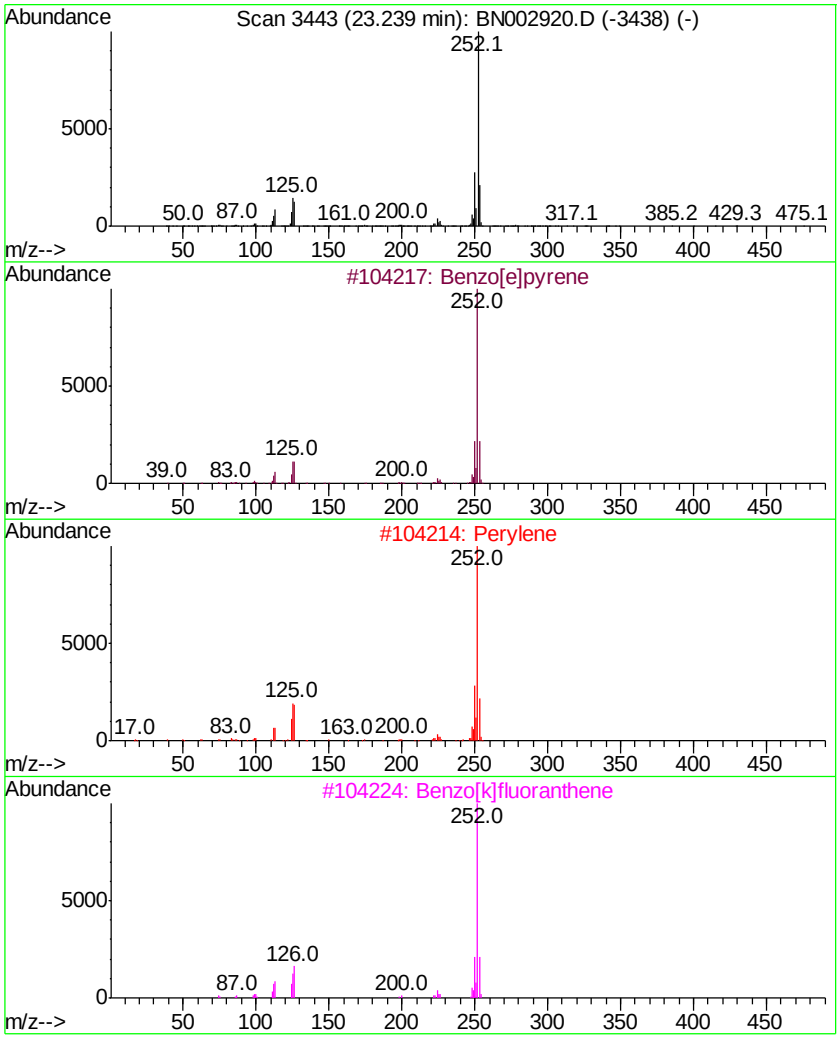
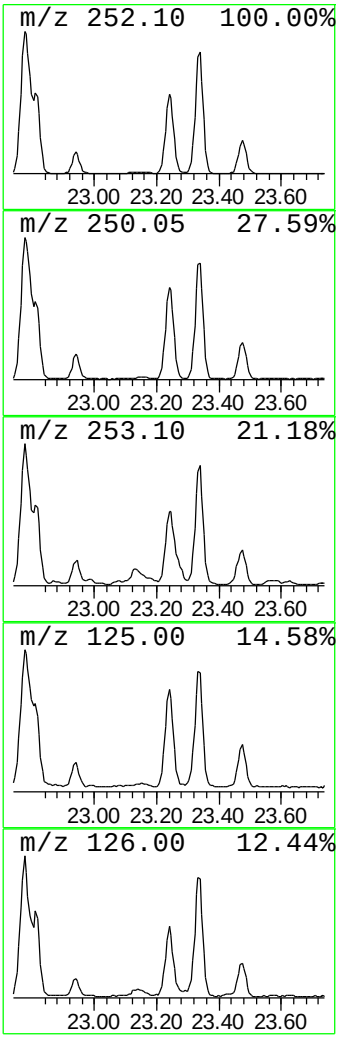
Peak Number 15 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	22.13 ng/ul	1721250	Perylene-d12	23.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Perylene	252	C20H12	000198-55-0	97
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
4	Benzo[a]pyrene	252	C20H12	000050-32-8	95
5	Perylene	252	C20H12	000198-55-0	93





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100418\
Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

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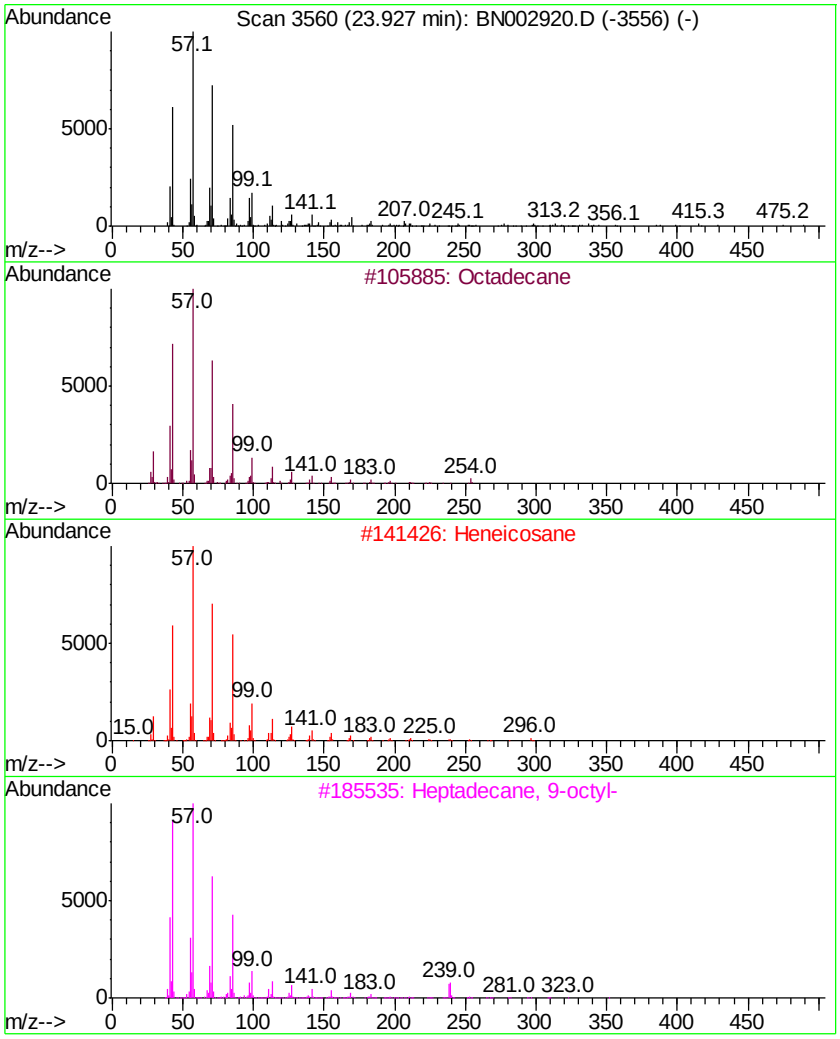
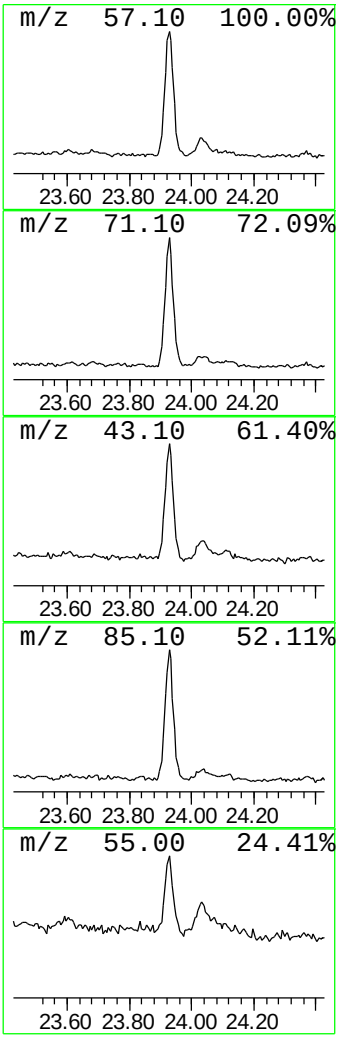
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.93	7.91 ng/ul	615411	Perylene-d12	23.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4	Docosane	310	C22H46	000629-97-0	87
5	Docosane	310	C22H46	000629-97-0	87





Tentatively Identified Compound (LSC) summary

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Data File : BN002920.D
Acq On : 04 Oct 2018 02:39
Operator : MJ/SJ
Sample : J5115-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : JLC63

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.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...	17.94	8.4	ng/ul	864843	4	17.02	2055650	20.0
4H-Cyclopenta[def...	18.09	7.7	ng/ul	788081	4	17.02	2055650	20.0
1,2-Benzenedicarb...	19.78	28.0	ng/ul	8351970	5	21.22	5965890	20.0
11H-Benzo[b]fluorene	19.95	3.0	ng/ul	885129	5	21.22	5965890	20.0
Cyclopenta(cd)pyr...	20.90	2.1	ng/ul	627194	5	21.22	5965890	20.0
Cyclopenta[cd]pyrene	20.94	4.5	ng/ul	1327160	5	21.22	5965890	20.0
Phosphoric acid, ...	21.72	2.2	ng/ul	664962	5	21.22	5965890	20.0
unknown-01	22.03	5.4	ng/ul	1602250	5	21.22	5965890	20.0
Acridine 10-oxide	22.05	5.1	ng/ul	1508560	5	21.22	5965890	20.0
unknown-02	22.21	5.2	ng/ul	1558320	5	21.22	5965890	20.0
unknown-03	22.39	10.9	ng/ul	848809	6	23.43	1555750	20.0
Benzo[e]pyrene	23.24	22.1	ng/ul	1721250	6	23.43	1555750	20.0
(DEL) Alkane: Str...	23.93	7.9	ng/ul	615411	6	23.43	1555750	20.0