LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_F\METHODS\8270-BF060517.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.557	9	12	51	rVB	1222824	2874345	100.00%	20.671%
2	4.457	501	505	517	rBV2	17949	50712	1.76%	0.365%
3	5.204	628	632	658	rBV	162370	536153	18.65%	3.856%
4	6.892	915	919	927	rVV	185490	445905	15.51%	3.207%
5	6.963	927	931	951	rVV	392880	913731	31.79%	6.571%
6	7.292	983	987	997	rBV	290003	410630	14.29%	2.953%
7	7.528	1023	1027	1043	rVB	408077	593505	20.65%	4.268%
8	8.216	1141	1144	1160	rBV	168763	374676	13.04%	2.695%
9	9.327	1328	1333	1345	rBV	344270	556398	19.36%	4.001%
10	9.427	1346	1350	1353	rVB2	24006	29125	1.01%	0.209%
11	10.410	1511	1517	1522	rBV	47698	54841	1.91%	0.394%
12	10.504	1527	1533	1540	rBV	25535	45006	1.57%	0.324%
13	11.104	1630	1635	1648	rBV	522372	679792	23.65%	4.889%
14	11.327	1669	1673	1677	rBV	33568	40016	1.39%	0.288%
15	11.810	1751	1755	1760	rBV	41437	66775	2.32%	0.480%
16	11.921	1770	1774	1782	rBV	45182	70316	2.45%	0.506%
17	12.145	1807	1812	1817	rBV2	376568	501914	17.46%	3.610%
18	12.186	1817	1819	1829	rVB3	35210	51983	1.81%	0.374%
19	13.451	2030	2034	2047	rBV	182528	309869	10.78%	2.228%
20	13.792	2085	2092	2097	rBV4	33682	78048	2.72%	0.561%
21	14.539	2215	2219	2223	rVV	387903	501698	17.45%	3.608%
22	14.580	2223	2226	2235	rVV	67536	106262	3.70%	0.764%
23	14.674	2238	2242	2252	rVB	54887	96224	3.35%	0.692%
24	15.398	2361	2365	2372	rVB3	25804	36252	1.26%	0.261%
25	15.562	2388	2393	2402	rVB5	22671	54322	1.89%	0.391%
26	15.804	2429	2434	2439	rBV4	21048	41506	1.44%	0.298%
27	15.886	2445	2448	2454	rBV4	15134	29317	1.02%	0.211%
28	16.162	2491	2495	2498	rBV	40968	50794	1.77%	0.365%
29	16.203	2499	2502	2507	rVB6	21150	37538	1.31%	0.270%
30	16.368	2524	2530	2537	rBV	213798	321901	11.20%	2.315%
31	16.509	2551	2554	2558	rVV3	23804	35446	1.23%	0.255%
32	16.668	2577	2581	2588	rVB	270162	345932	12.04%	2.488%
33	16.727	2588	2591	2594	rBV5	20384	29016	1.01%	0.209%
34	16.921	2619	2624	2633	rVV	471901	604833	21.04%	4.350%

LSC Area Percent Report

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Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_F\METHODS\8270-BF060517.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.998	2635	2637	2641	rVB3	30427	35422	1.23%	0.255%
36	17.068	2646	2649	2652	rBV4	25118	36106	1.26%	0.260%
37	17.121	2653	2658	2659	rBV4	28505	46884	1.63%	0.337%
38	17.286	2681	2686	2689	rBV2	66627	118390	4.12%	0.851%
39	17.327	2691	2693	2697	rVV4	24206	37587	1.31%	0.270%
40	17.786	2768	2771	2774	rVB5	34498	44522	1.55%	0.320%
41	17.839	2774	2780	2784	rBV2	63197	97996	3.41%	0.705%
42	17.945	2795	2798	2799	rBV2	52590	46147	1.61%	0.332%
43	17.980	2802	2804	2805	rVB2	38424	29281	1.02%	0.211%
44	18.121	2825	2828	2831	rBV2	88642	120239	4.18%	0.865%
45	18.186	2836	2839	2840	rBV3	45002	52840	1.84%	0.380%
46	18.268	2848	2853	2858	rBV2	505099	839164	29.19%	6.035%
47	18.668	2917	2921	2925	rVB6	64668	115041	4.00%	0.827%
48	18.733	2925	2932	2933	rBV6	61695	124453	4.33%	0.895%
49	19.303	3027	3029	3032	rVB2	90409	79202	2.76%	0.570%
50	19.397	3042	3045	3047	rBV4	79659	104266	3.63%	0.750%
51	19.544	3067	3070	3073	rBV	222301	198768	6.92%	1.429%
52	19.839	3116	3120	3122	rBV	191649	241871	8.41%	1.739%
53	19.897	3128	3130	3133	rVB	178741	194377	6.76%	1.398%
54	19.944	3135	3138	3142	rBV2	273156	367644	12.79%	2.644%

Sum of corrected areas: 13904981

LSC Report - Integrated Chromatogram

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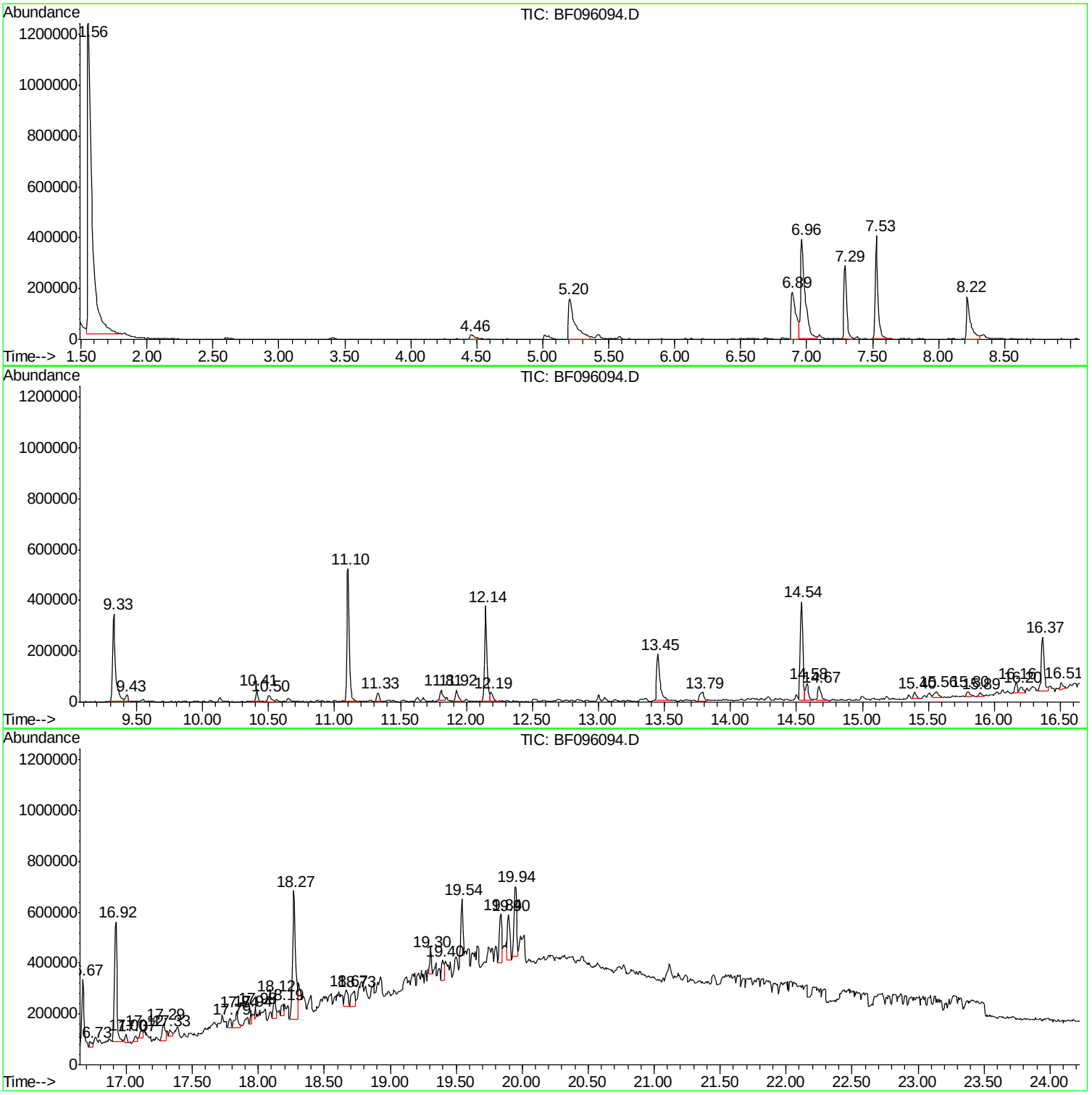
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

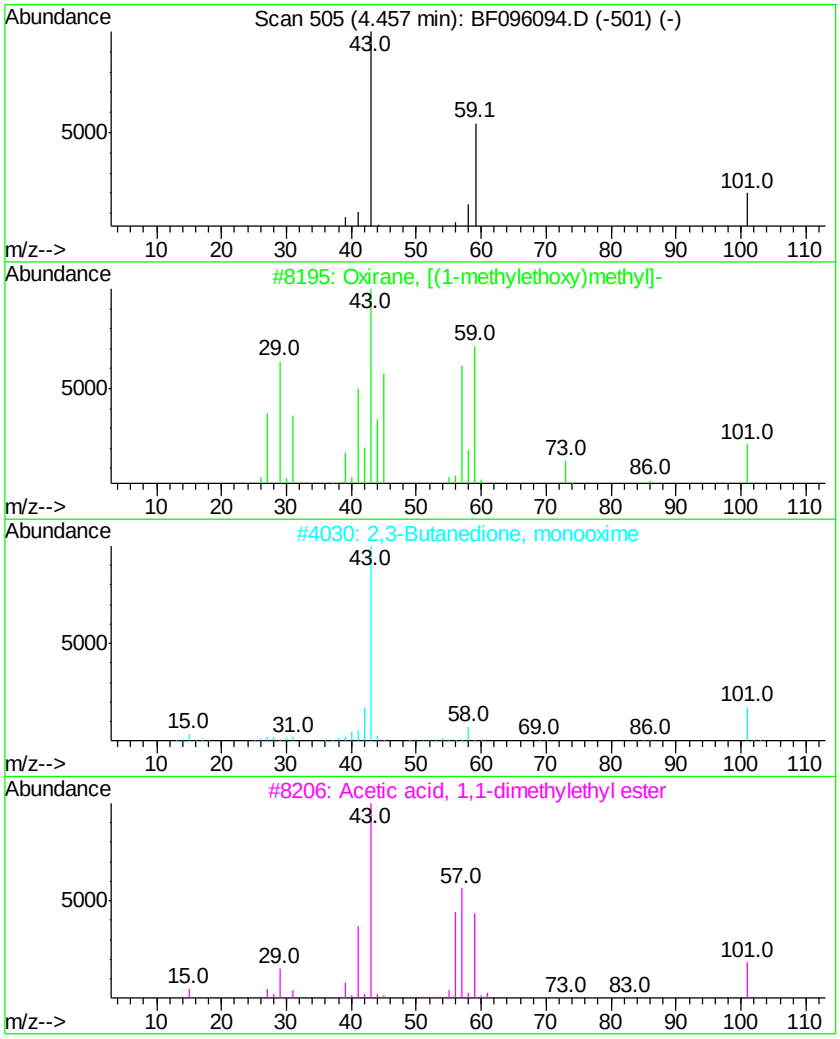
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

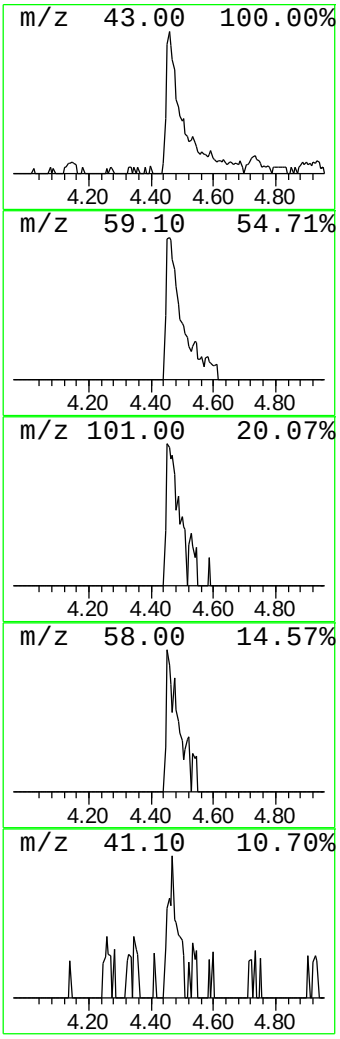
****

Peak Number 2 Oxirane, [(1-methylethoxy)m... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	2.47 ng	50712	1,4-Dichlorobenzene-d4	7.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	64
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	9
5	3-Hexanol	102	C6H14O	000623-37-0	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

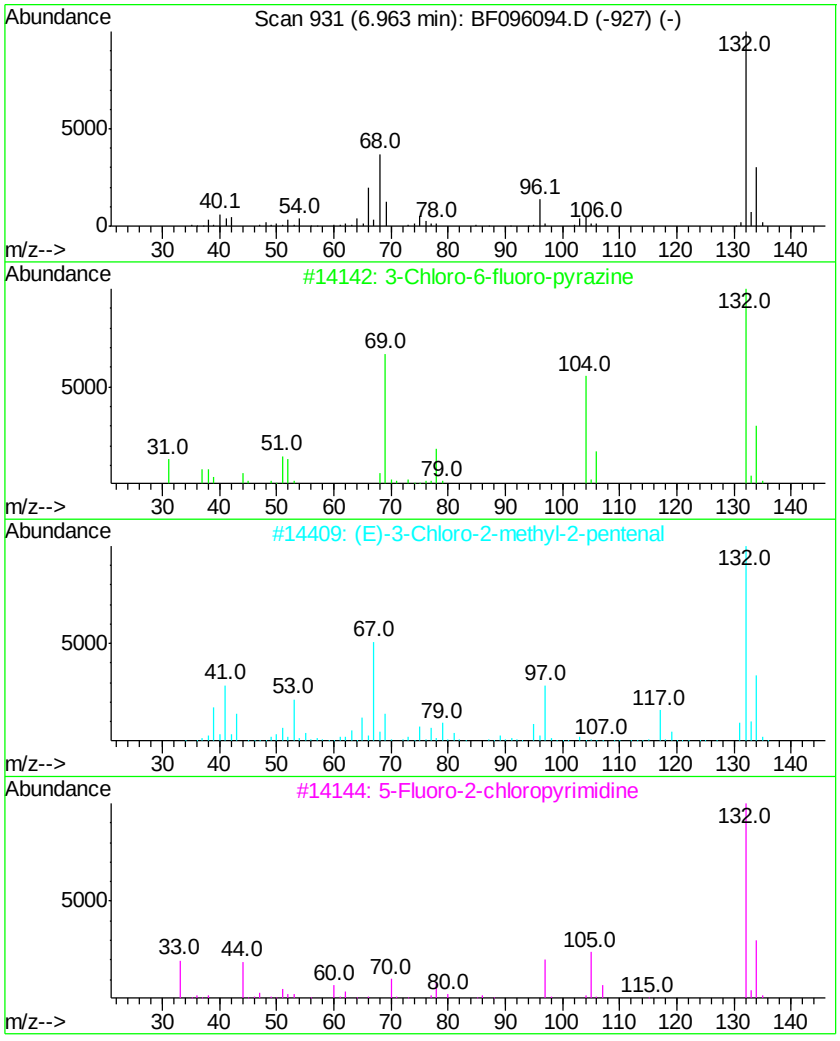
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

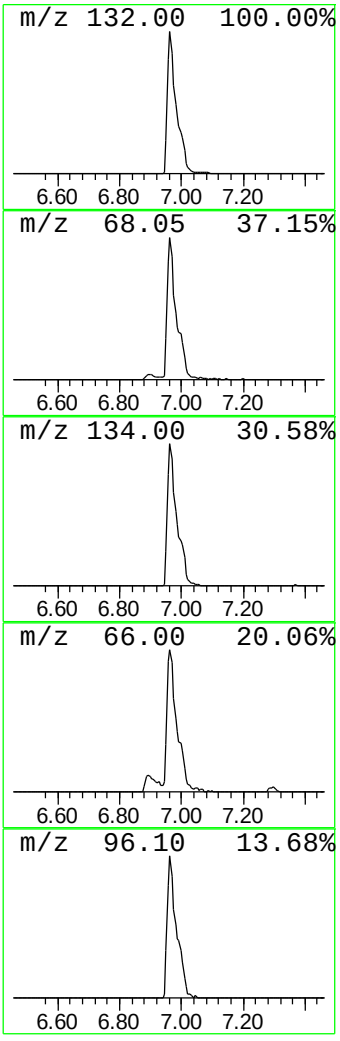
****

Peak Number 3 unknown6.96 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.96	44.50 ng	913731	1,4-Dichlorobenzene-d4	7.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
4	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5	Xenon	132	Xe	007440-63-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
Data File : BF096094.D
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Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

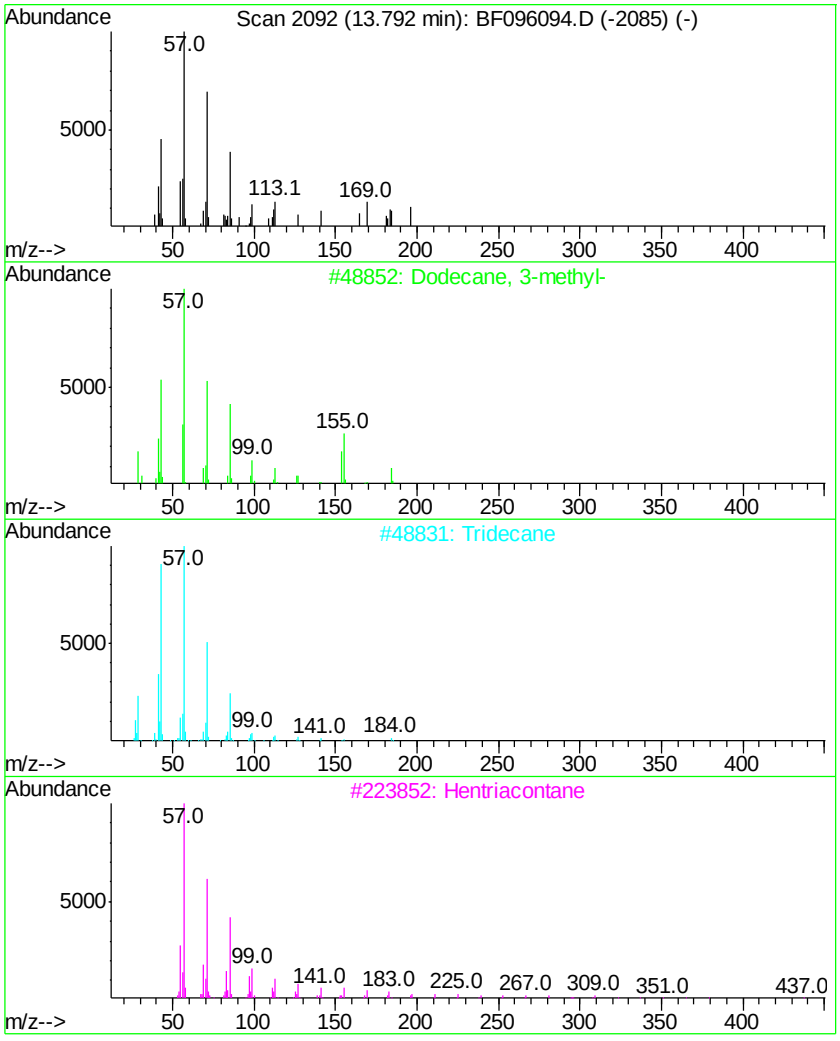
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

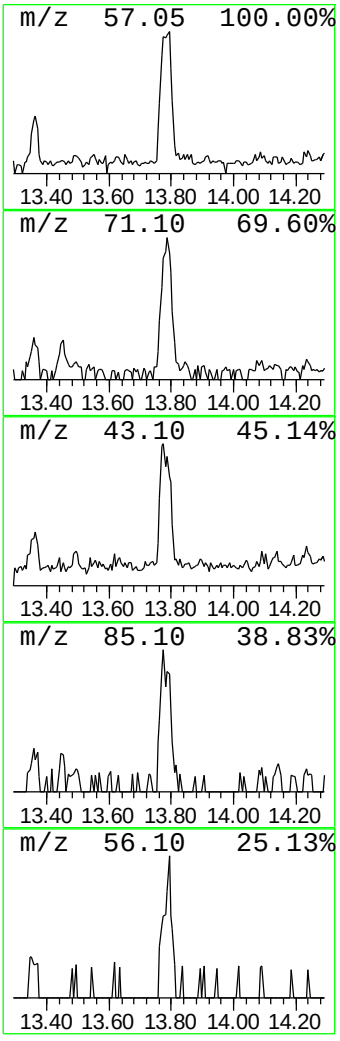
****

Peak Number 5 Dodecane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.11 ng	78048	Phenanthrene-d10	14.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 3-methyl-	184	C13H28	017312-57-1	76
2	Tridecane	184	C13H28	000629-50-5	68
3	Hentriacontane	437	C31H64	000630-04-6	64
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	58
5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	58





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

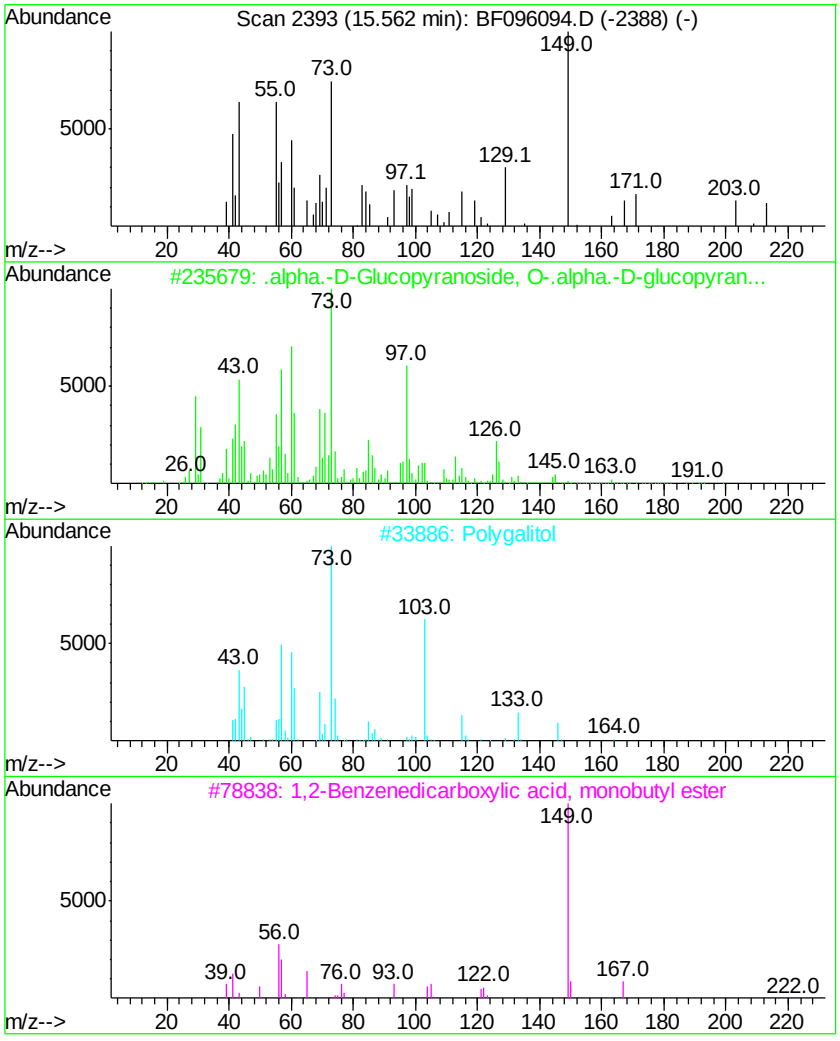
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

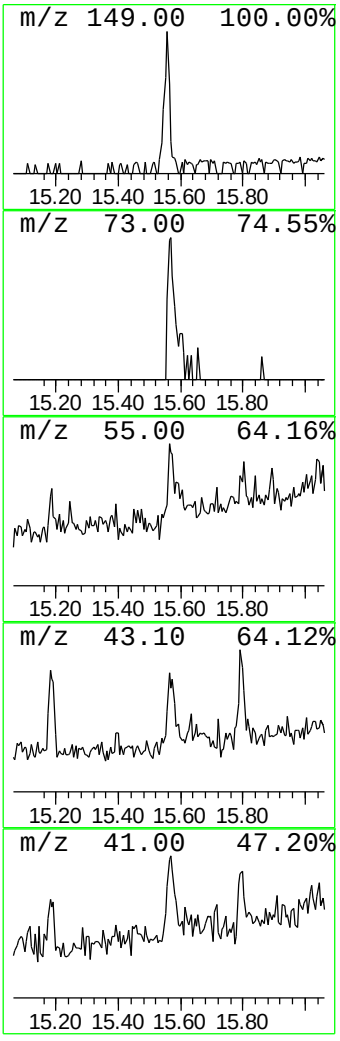
****

Peak Number 6 unknown15.56 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	2.17 ng	54322	Phenanthrene-d10	14.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-D-Glucopyranoside, O-.al...	504	C18H32O16	000597-12-6	16
2	Polygalitol	164	C6H12O5	1000126-63-2	14
3	1,2-Benzenedicarboxylic acid, mo...	222	C12H14O4	000131-70-4	14
4	Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	12
5	Methyl 4,6-ethylidene-.alpha.-d-...	220	C9H16O6	1000102-53-5	12





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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

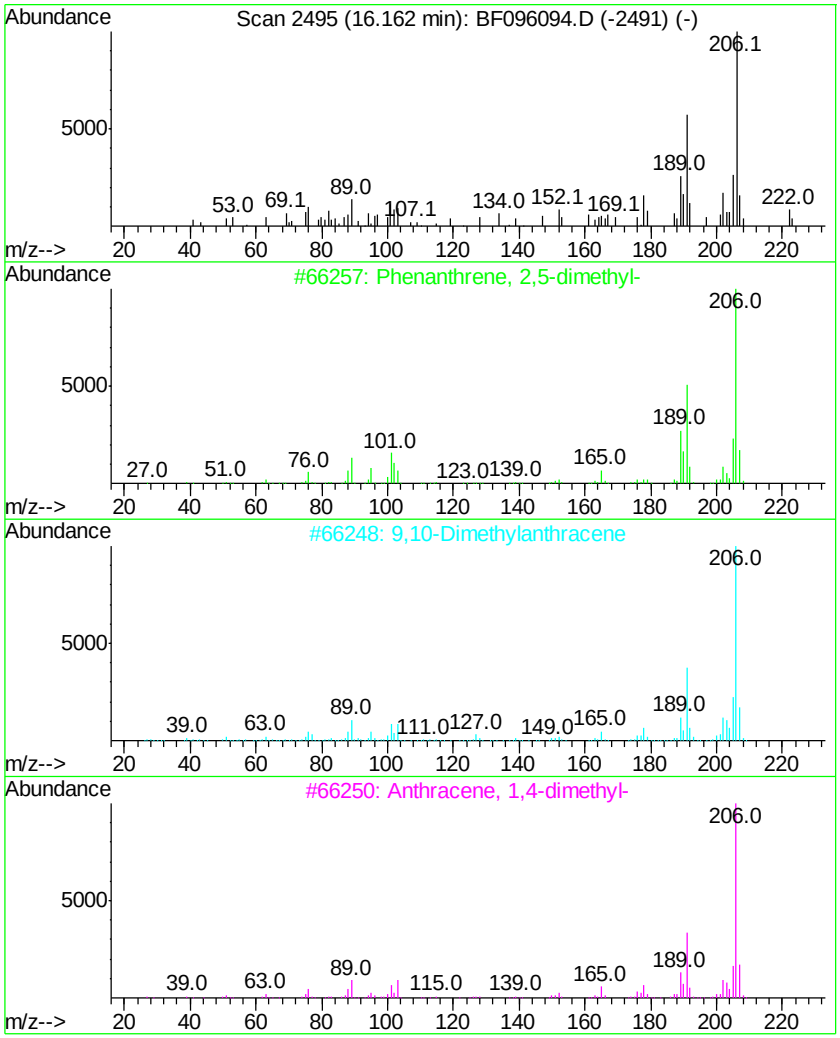
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

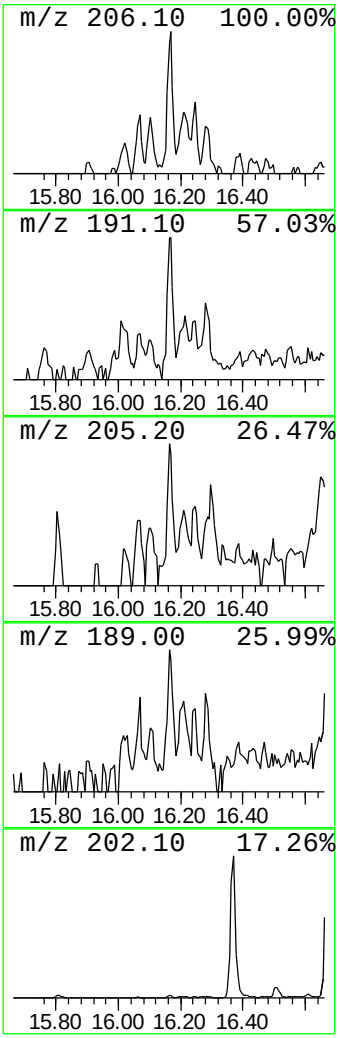
****

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	2.02 ng	50794	Phenanthrene-d10	14.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
3	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
4	di-p-Tolylacetylene	206	C16H14	002789-88-0	90
5	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	89





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ClientSampleId : G21-0-1.5

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TIC Library : C:\DATABASE\NIST11.L
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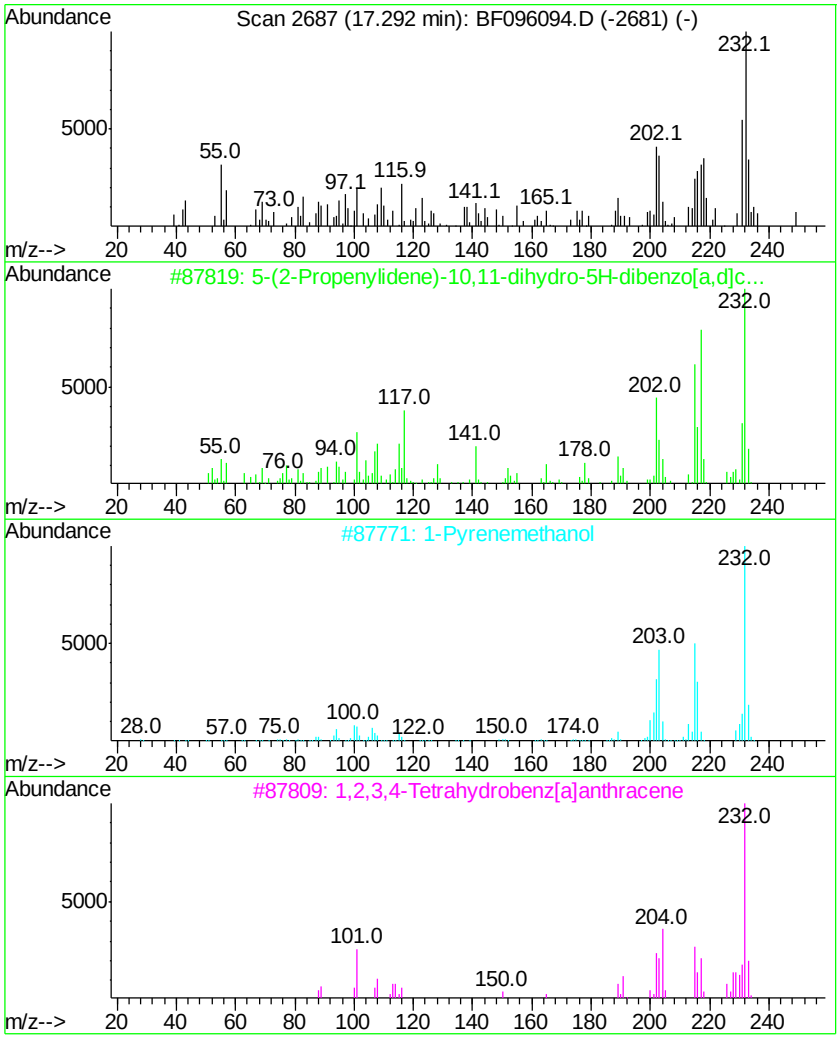
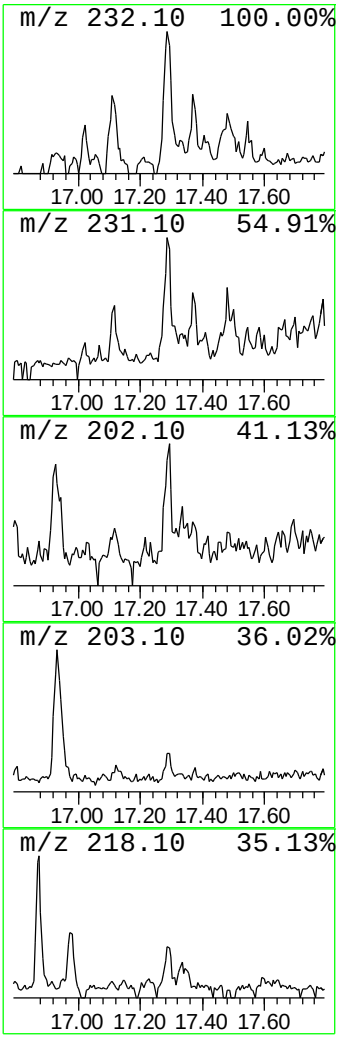
Peak Number 8 unknown17.29 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.29	2.82 ng	118390	Chrysene-d12	18.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(2-Propenylidene)-10,11-dihydr...	232	C18H16	024755-73-5	49
2	1-Pyrenemethanol	232	C17H12O	024463-15-8	46
3	1,2,3,4-Tetrahydrobenz[a]anthracene	232	C18H16	004483-98-1	43
4	Benzoic acid, 3-(5-formyl-2-furyl)-	216	C12H8O4	304884-54-6	42
5	trans-1,1'-Bibenzoindanylidene	232	C18H16	024536-68-3	41





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Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

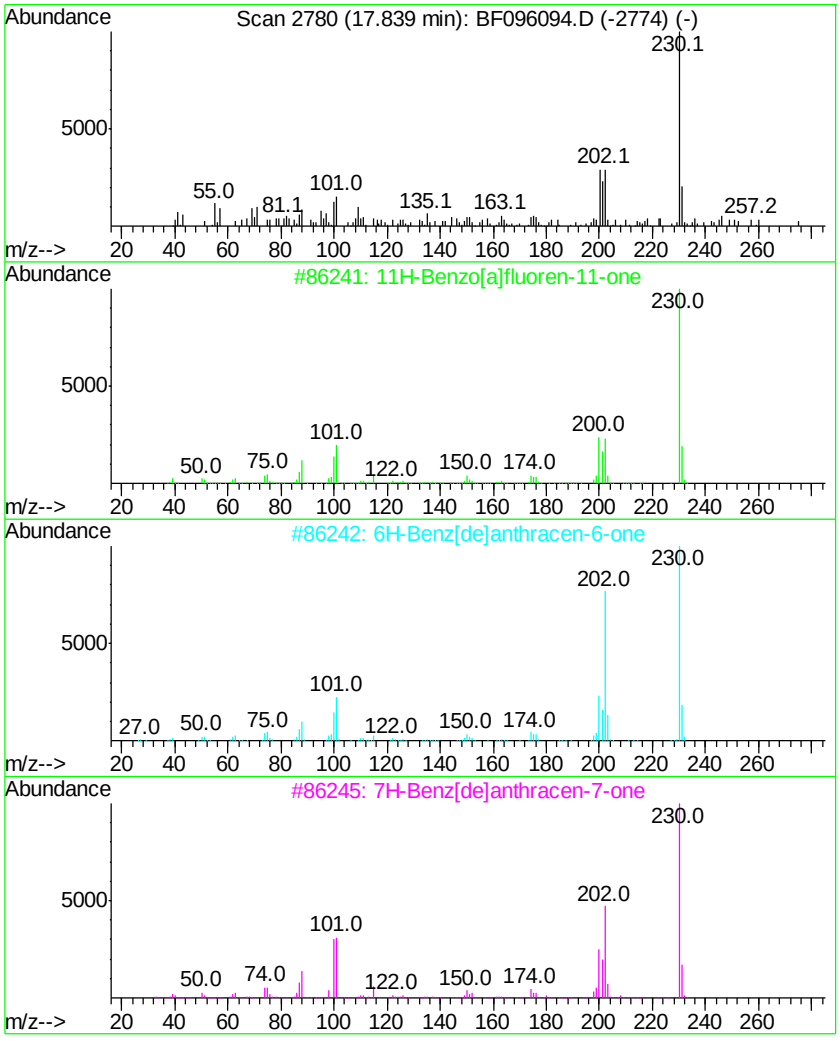
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

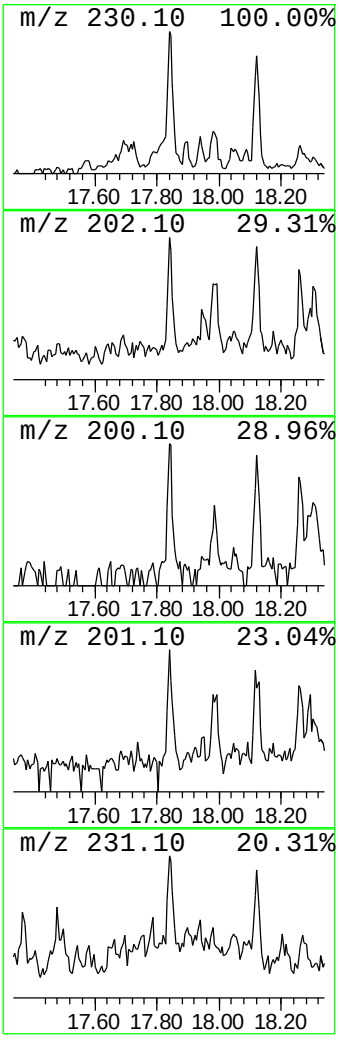
****

Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.84	2.34 ng	97996	Chrysene-d12	18.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2	6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	86
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	64
4	Dibenzo[c,h][2,6]naphthyridine	230	C16H10N2	000218-30-4	50
5	o-Terphenyl	230	C18H14	000084-15-1	47





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Acq On : 21 Jun 2017 20:23
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ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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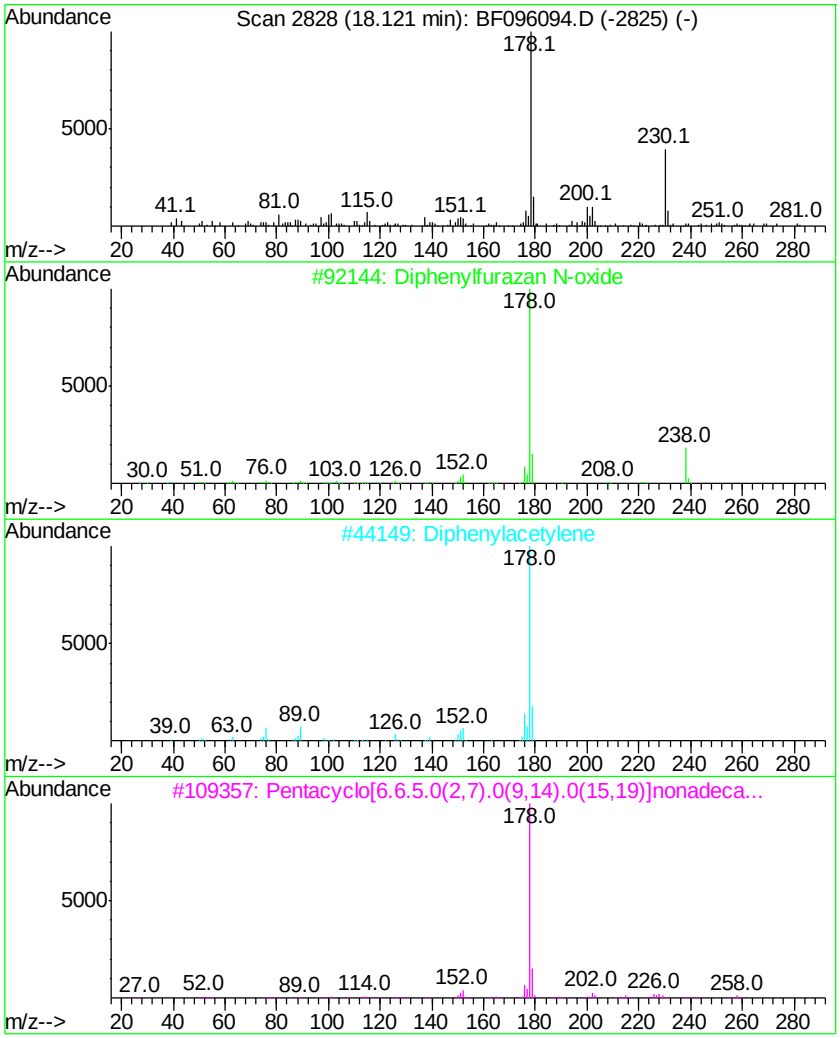
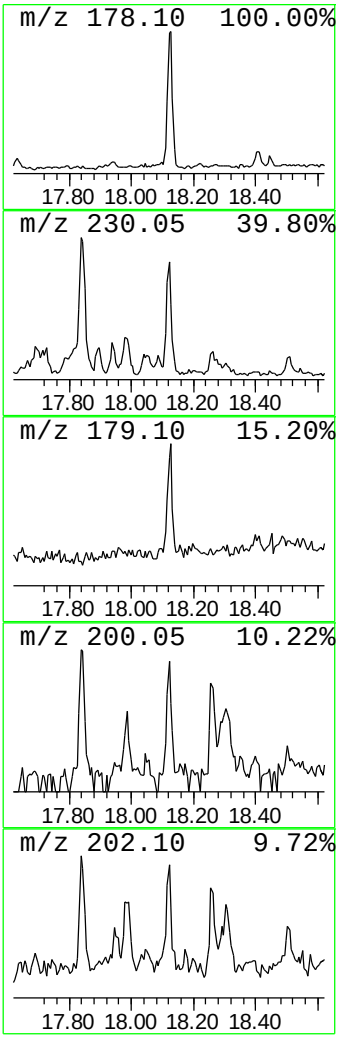
****

Peak Number 10 Diphenylfurazan N-oxide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.12	2.87 ng	120239	Chrysene-d12	18.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenylfurazan N-oxide	238	C14H10N2O2	005585-14-8	55
2	Diphenylacetylene	178	C14H10	000501-65-5	55
3	Pentacyclo[6.6.5.0(2,7).0(9,14)....	258	C19H14O	1000156-78-8	49
4	Anthracene	178	C14H10	000120-12-7	49
5	2,3-Naphthalenedicarbonitrile	178	C12H6N2	022856-30-0	46





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Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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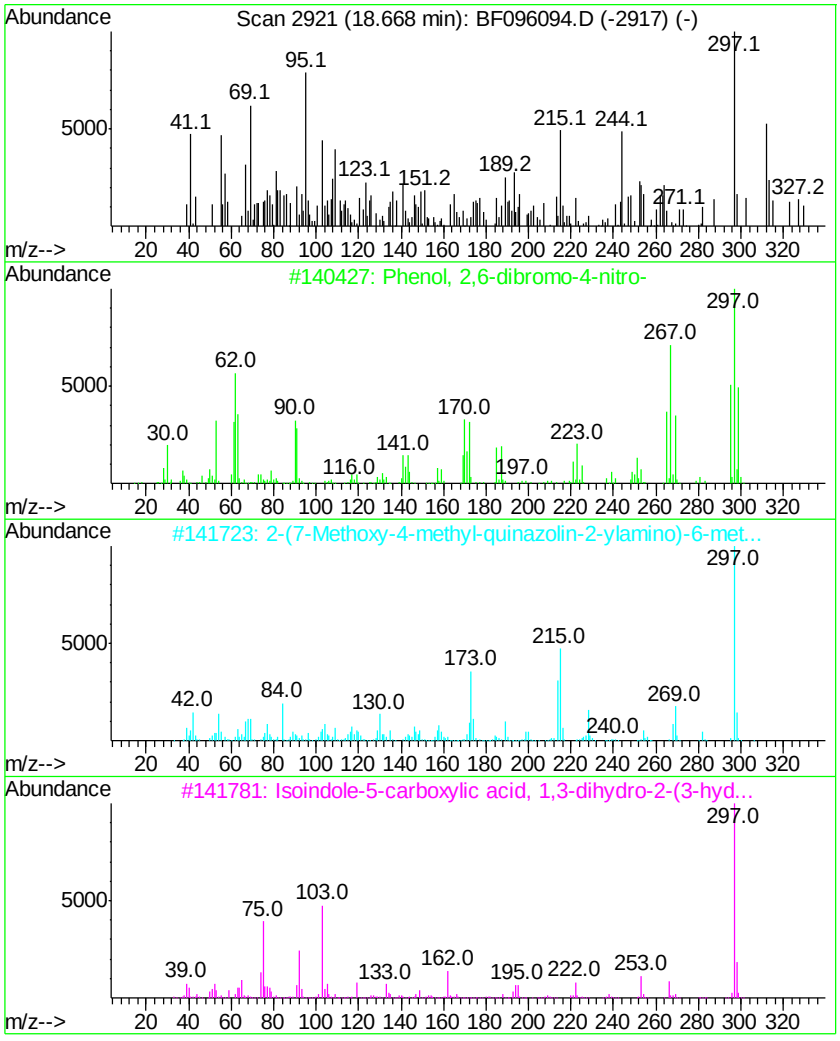
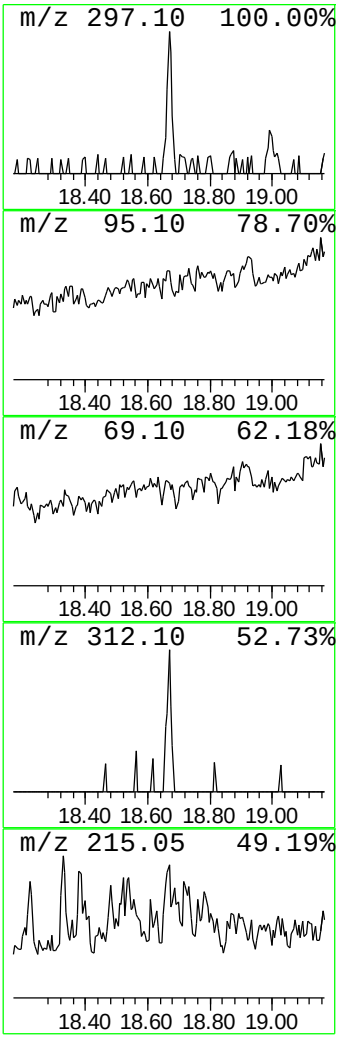
Peak Number 11 unknown18.67 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.67	2.74 ng	115041	Chrysene-d12	18.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dibromo-4-nitro-	295	C6H3Br2NO3	000099-28-5	25
2	2-(7-Methoxy-4-methyl-quinazolin...	297	C15H15N5O2	1000301-28-4	18
3	Isoindole-5-carboxylic acid, 1,3...	297	C16H11NO5	339339-09-2	18
4	4'-Hydroxy-3'-methoxyacetophenon...	312	C12H9F5O4	1000375-93-3	14
5	Butylphosphonic acid, butyl 4-is...	312	C17H29O3P	1000322-96-4	14





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Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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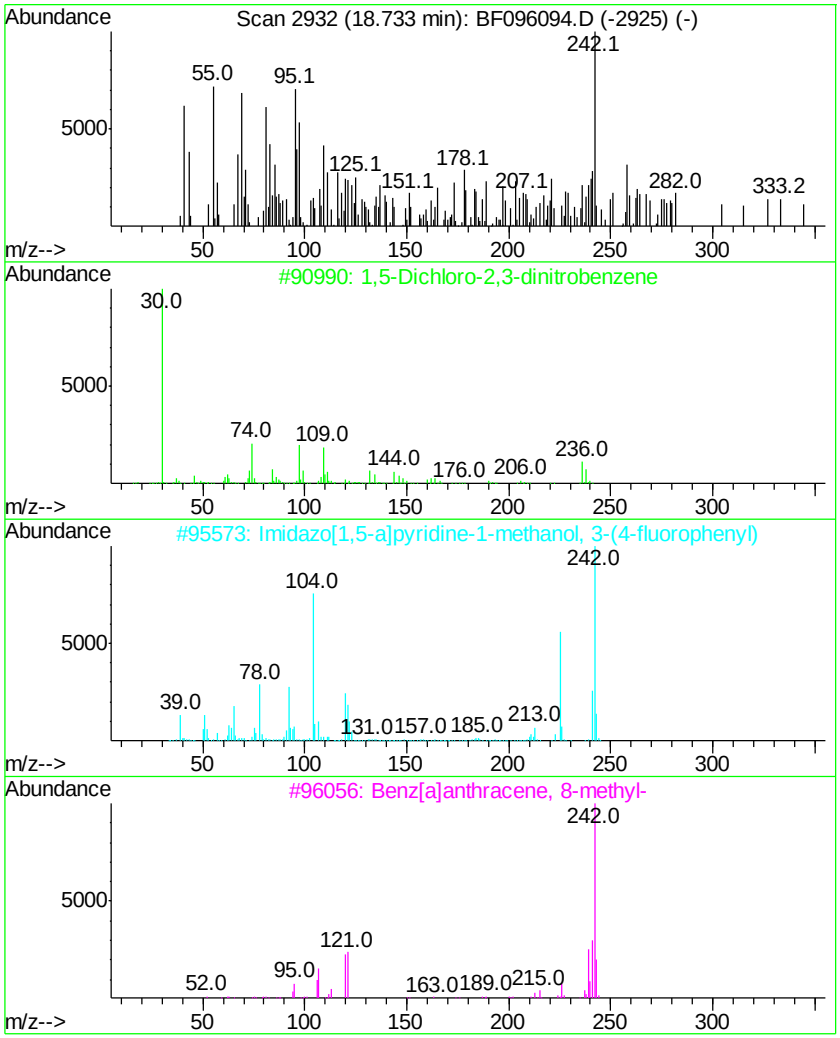
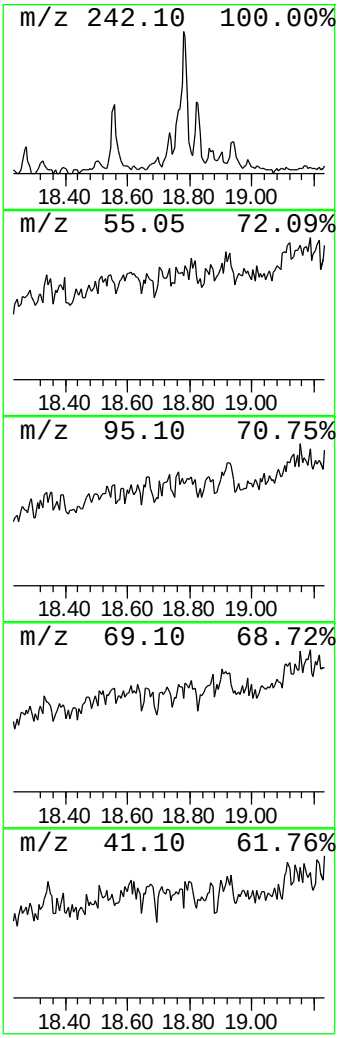
Peak Number 12 unknown18.73 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	2.97 ng	124453	Chrysene-d12	18.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Dichloro-2,3-dinitrobenzene	236	C6H2Cl2N2O4	028689-08-9	18
2	Imidazo[1,5-a]pyridine-1-methano...	242	C14H11FN2O	1000337-93-8	15
3	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	15
4	Octadec-9-enoic acid	282	C18H34O2	1000190-13-7	14
5	Pyridazino[4,5-g]phtalazine-1,6(...	242	C12H10N4O2	312520-09-5	11





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Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
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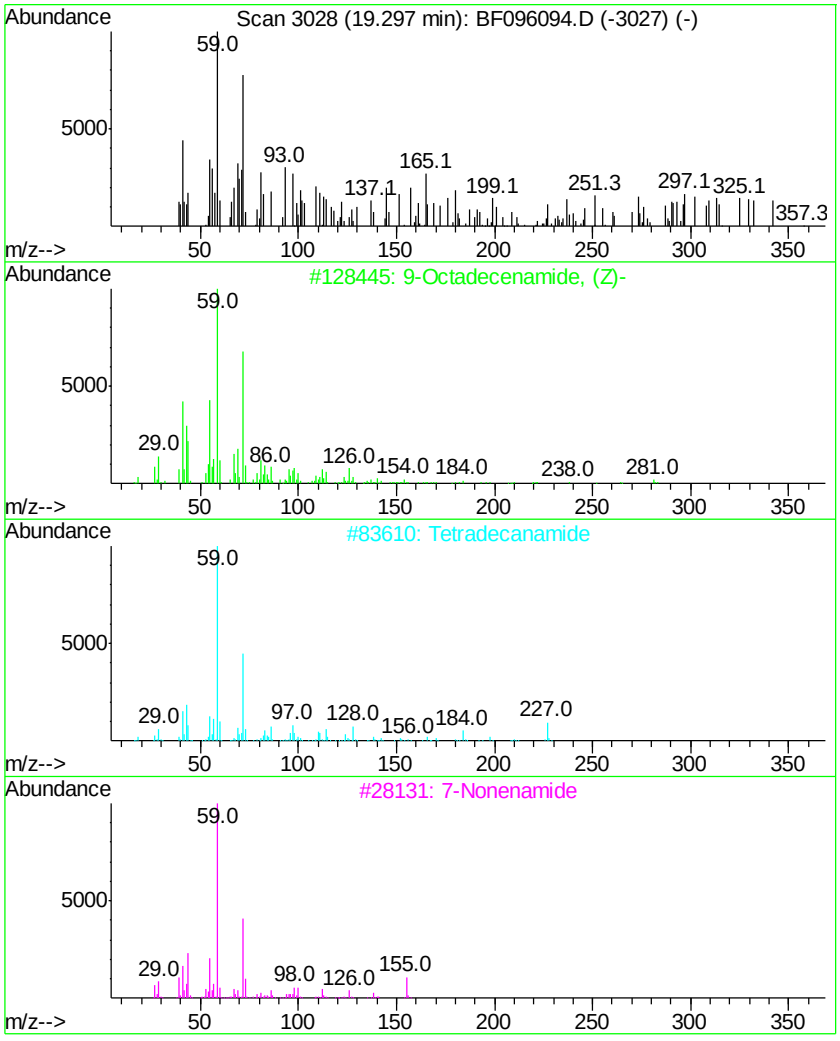
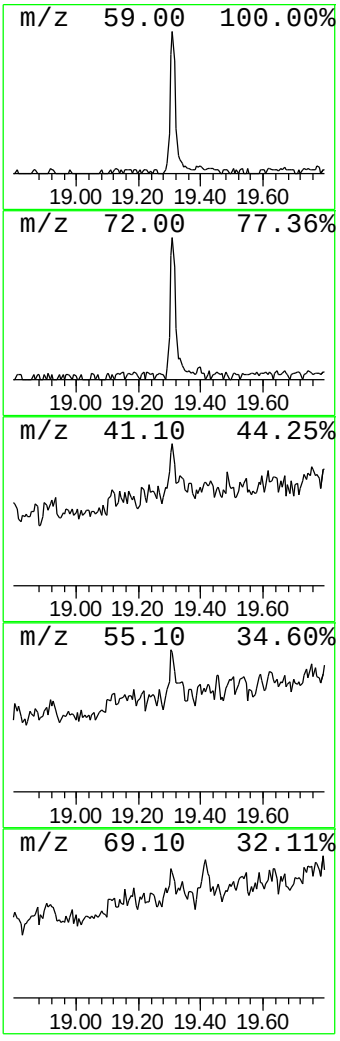
Peak Number 13 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.30	4.31 ng	79202	Perylene-d12	19.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
2	Tetradecanamide	227	C14H29NO	000638-58-4	47
3	7-Nonenamide	155	C9H17NO	090949-53-4	46
4	Nonanamide	157	C9H19NO	001120-07-6	43
5	Octanamide	143	C8H17NO	000629-01-6	38





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Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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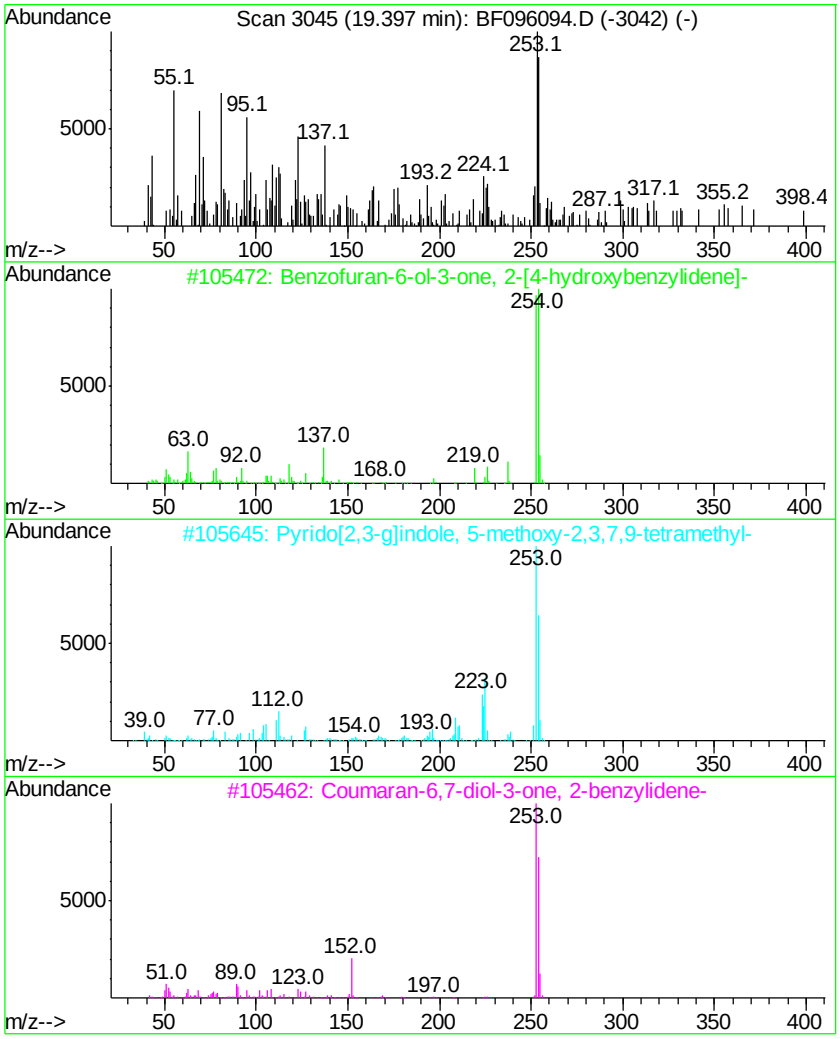
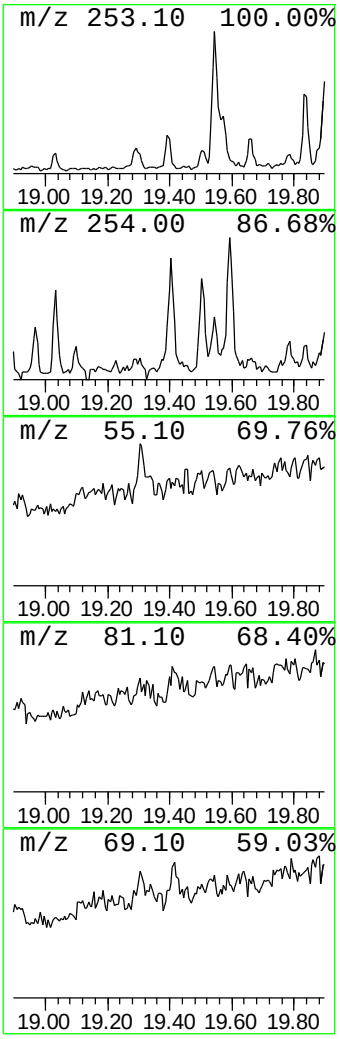
Peak Number 14 unknown19.40 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.40	5.67 ng	104266	Perylene-d12	19.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran-6-ol-3-one, 2-[4-hydr...	254	C15H10O4	1000128-64-9	43
2	Pyrido[2,3-g]indole, 5-methoxy-2...	254	C16H18N2O	201991-45-9	43
3	Coumaran-6,7-diol-3-one, 2-benzy...	254	C15H10O4	029813-90-9	38
4	1,1'-Binaphthalene	254	C20H14	000604-53-5	38
5	1H-Pyrrolo[2,3-f]quinolin-9-ol, ...	254	C16H18N2O	1000303-71-9	35





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Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
Sample : I3736-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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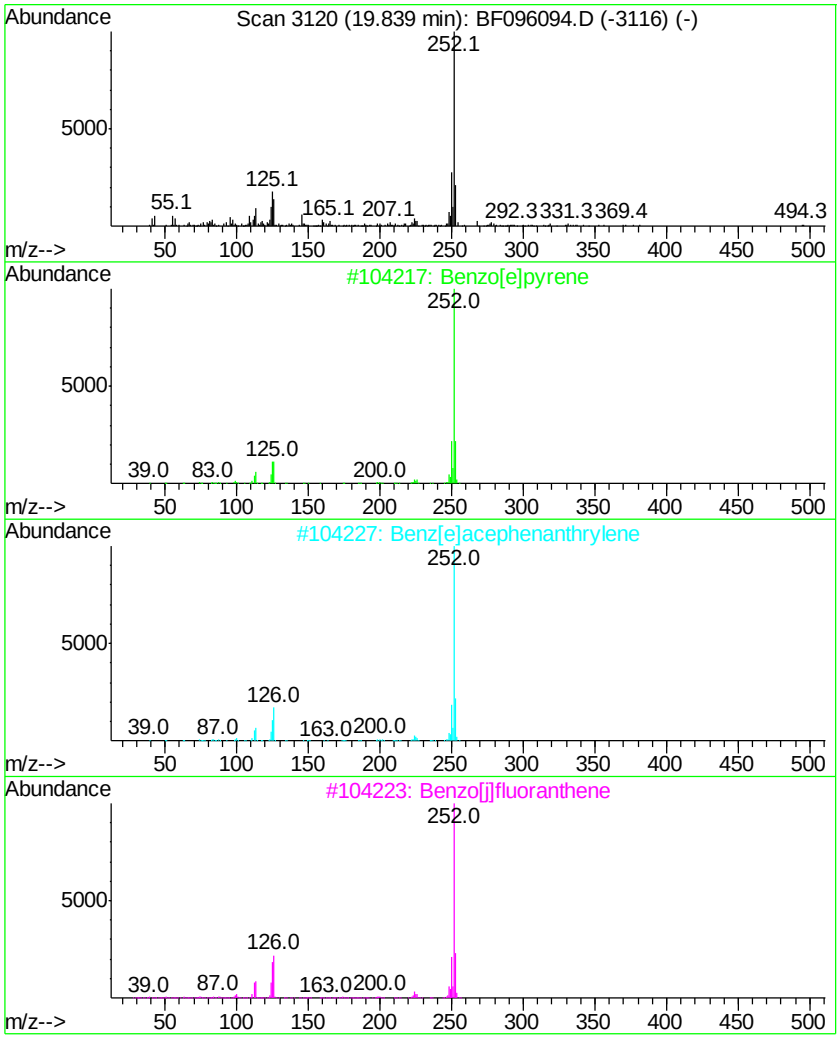
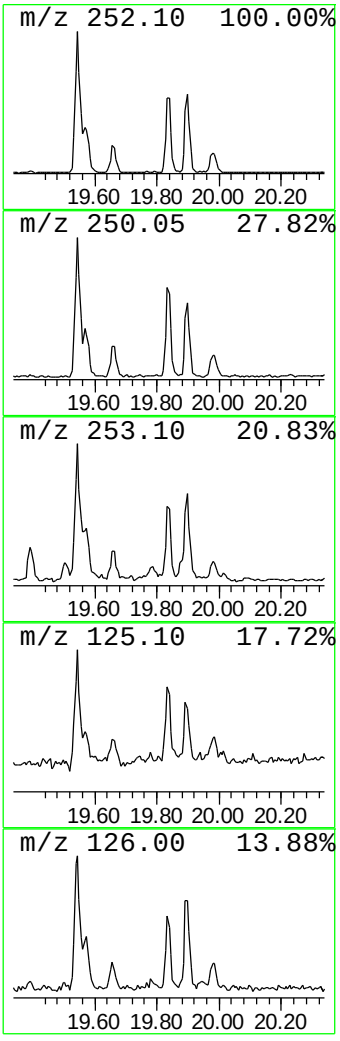
Peak Number 15 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.84	13.16 ng	241871	Perylene-d12	19.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	93
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





Tentatively Identified Compound (LSC) summary

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Data File : BF096094.D
Acq On : 21 Jun 2017 20:23
Operator : SJ/MA
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G21-0-1.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Oxirane, [(1-meth...	4.46	2.5	ng	50712	1	7.29	410630	20.0
unknown6.96	6.96	44.5	ng	913731	1	7.29	410630	20.0
Dodecane, 3-methyl-	13.79	3.1	ng	78048	4	14.54	501698	20.0
unknown15.56	15.56	2.2	ng	54322	4	14.54	501698	20.0
Phenanthrene, 2,5...	16.16	2.0	ng	50794	4	14.54	501698	20.0
unknown17.29	17.29	2.8	ng	118390	5	18.27	839164	20.0
11H-Benzo[a]fluor...	17.84	2.3	ng	97996	5	18.27	839164	20.0
Diphenylfurazan N...	18.12	2.9	ng	120239	5	18.27	839164	20.0
unknown18.67	18.67	2.7	ng	115041	5	18.27	839164	20.0
unknown18.73	18.73	3.0	ng	124453	5	18.27	839164	20.0
9-Octadecenamide,...	19.30	4.3	ng	79202	6	19.95	367644	20.0
unknown19.40	19.40	5.7	ng	104266	6	19.95	367644	20.0
Benzo[e]pyrene	19.84	13.2	ng	241871	6	19.95	367644	20.0