

Data Path : Z:\HPCHEM1\BNA F\DATA\BF043018\
 Data File : BF104947.D
 Acq On : 30 Apr 2018 21:41
 Operator : JU/SJ
 Sample : J2609-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01A

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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	5.416	562	566	587	rBV	2042747	2162328	15.01%	4.913%
2	6.439	736	740	757	rBV	2002998	2202938	15.29%	5.005%
3	6.563	757	761	765	rBV	2279088	2219603	15.40%	5.043%
4	6.786	796	799	803	rBV	772914	645078	4.48%	1.466%
5	6.945	822	826	829	rBV	2057118	1783893	12.38%	4.053%
6	7.357	889	896	899	rBV	1478604	1353818	9.40%	3.076%
7	8.075	1014	1018	1021	rBV	958907	842939	5.85%	1.915%
8	9.151	1196	1201	1204	rBV	2582456	2407489	16.71%	5.470%
9	9.545	1264	1268	1275	rVB	228002	216411	1.50%	0.492%
10	9.827	1312	1316	1319	rBV2	1043312	924099	6.41%	2.099%
11	10.621	1447	1451	1455	rBV	1444798	1381997	9.59%	3.140%
12	10.721	1466	1468	1478	rVB	94341	157704	1.09%	0.358%
13	11.321	1566	1570	1572	rBV	920114	875772	6.08%	1.990%
14	11.345	1572	1574	1578	rVB	991446	752120	5.22%	1.709%
15	11.398	1580	1583	1585	rBV	161655	154590	1.07%	0.351%
16	11.663	1624	1628	1632	rVB2	131138	151752	1.05%	0.345%
17	11.845	1656	1659	1662	rVV2	259017	270555	1.88%	0.615%
18	11.933	1671	1674	1677	rBV	257270	245249	1.70%	0.557%
19	12.539	1773	1777	1780	rBV	1139201	968701	6.72%	2.201%
20	12.627	1789	1792	1796	rVB2	174175	158442	1.10%	0.360%
21	12.768	1810	1816	1819	rBV	877798	921242	6.39%	2.093%
22	12.910	1836	1840	1846	rVB	1768342	1775611	12.32%	4.034%
23	13.092	1867	1871	1874	rVB2	171651	242030	1.68%	0.550%
24	13.715	1974	1977	1980	rBV2	144979	157089	1.09%	0.357%
25	13.780	1986	1988	1993	rBV2	124427	188594	1.31%	0.428%
26	13.974	2011	2021	2023	rBV	5474504	14409957	100.00%	32.738%
27	13.998	2023	2025	2030	rVB	473598	447933	3.11%	1.018%
28	14.221	2060	2063	2067	rBV3	124058	179558	1.25%	0.408%
29	14.562	2118	2121	2123	rBV	280834	273722	1.90%	0.622%
30	14.592	2123	2126	2130	rVB	1017975	893653	6.20%	2.030%
31	14.921	2179	2182	2185	rBV	210932	242552	1.68%	0.551%
32	14.956	2185	2188	2190	rBV2	152185	156115	1.08%	0.355%
33	15.015	2195	2198	2201	rBV	258879	341403	2.37%	0.776%
34	15.309	2244	2248	2256	rVB3	256725	518934	3.60%	1.179%

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Integration Parameters: rteint.p
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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-BF040618.M
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35	15.380	2256	2260	2265	rBV	219579	270881	1.88%	0.615%
36	15.439	2266	2270	2274	rBV	393890	465043	3.23%	1.057%
37	16.039	2368	2372	2378	rVB	159862	240324	1.67%	0.546%
38	16.898	2513	2518	2523	rBV	106207	200073	1.39%	0.455%
39	17.833	2667	2677	2689	rVB	836795	2215815	15.38%	5.034%

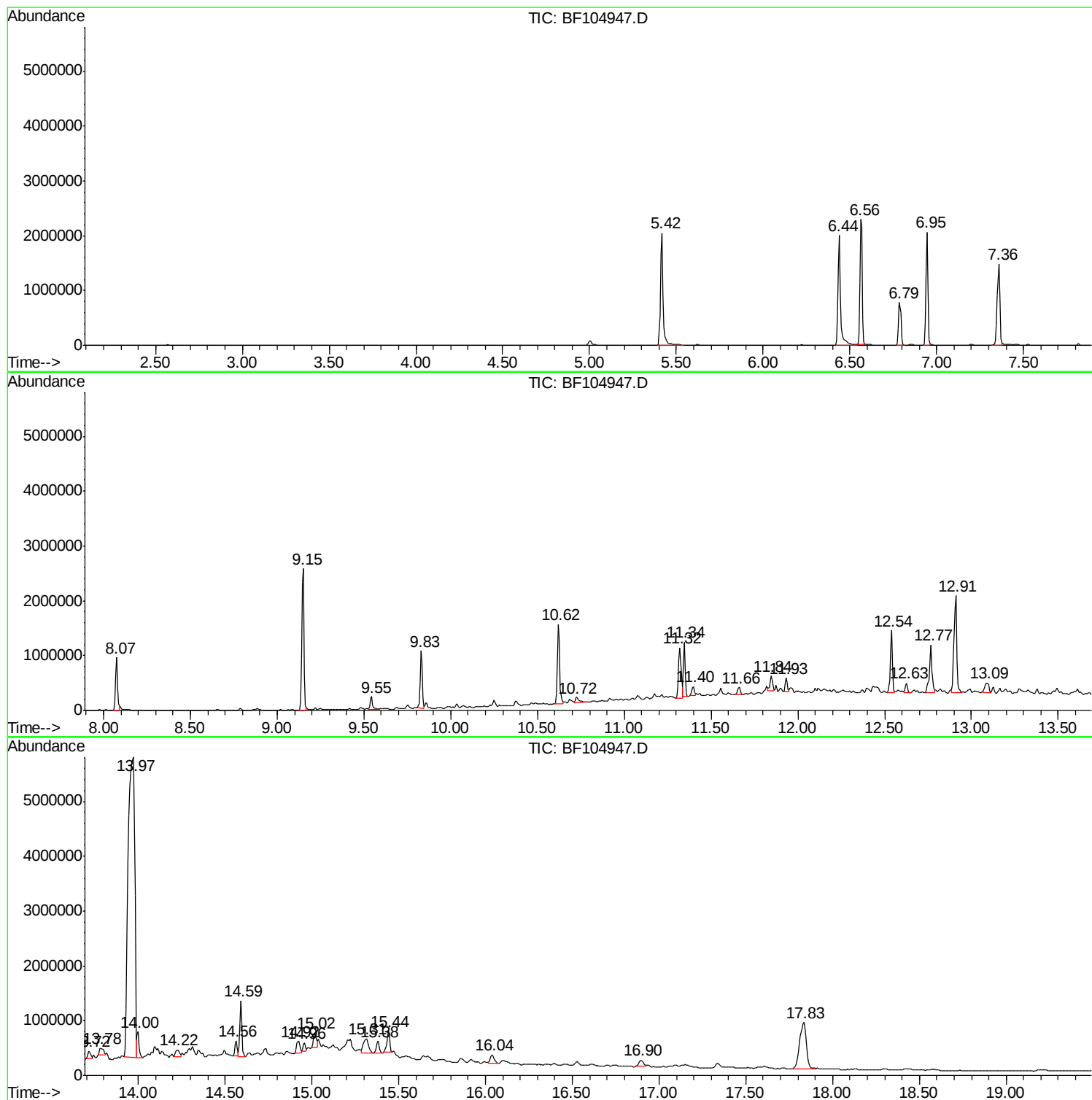
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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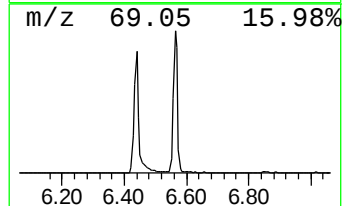
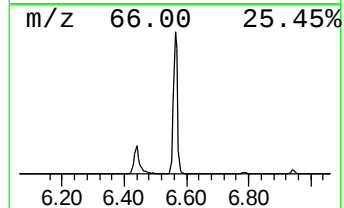
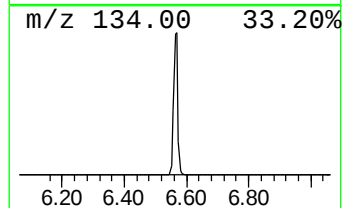
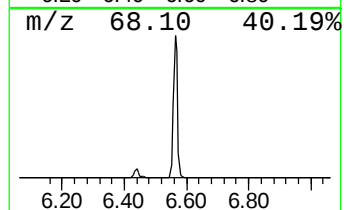
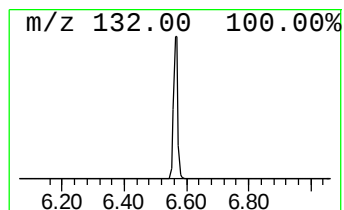
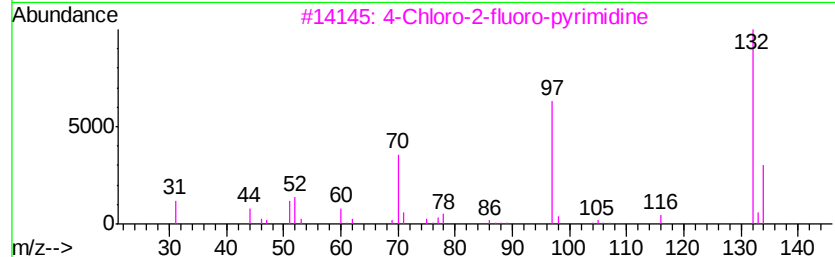
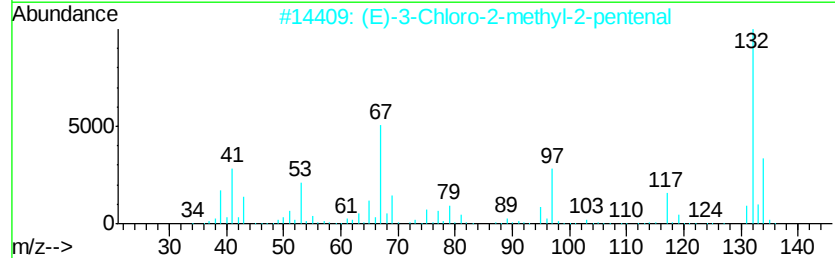
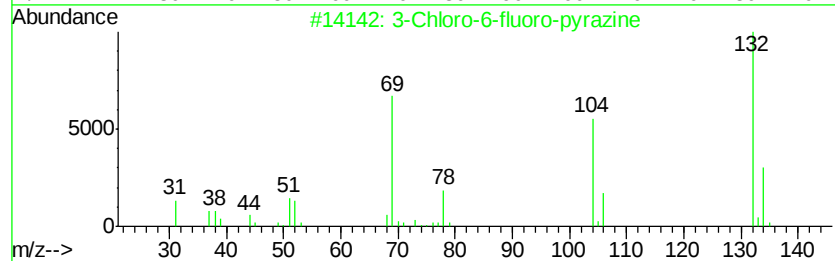
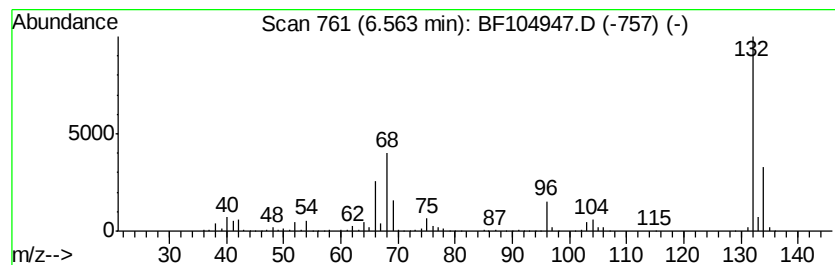
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	68.82 ng	2219600	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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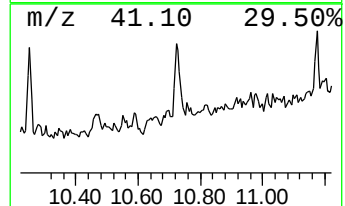
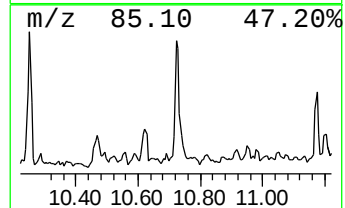
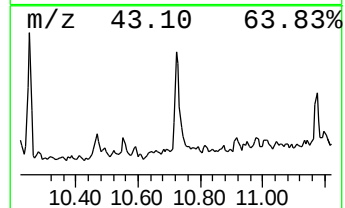
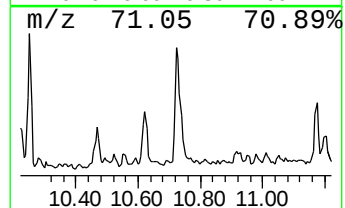
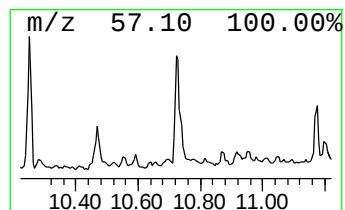
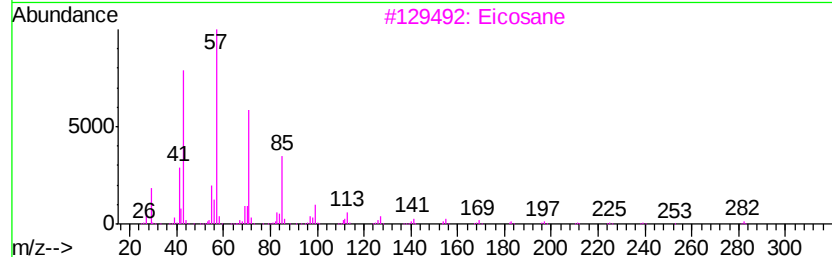
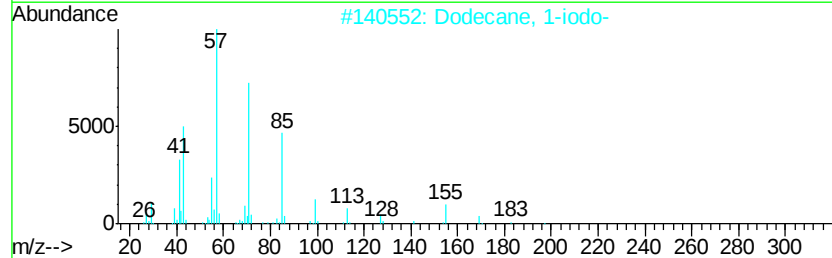
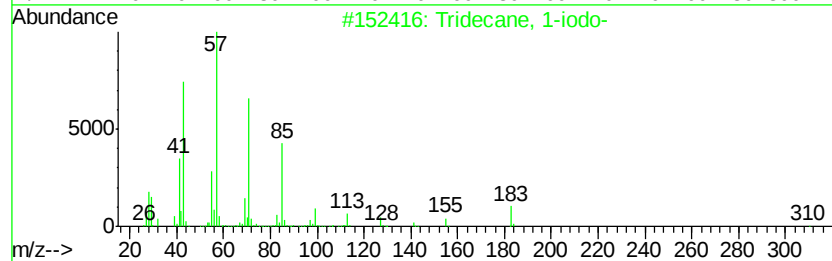
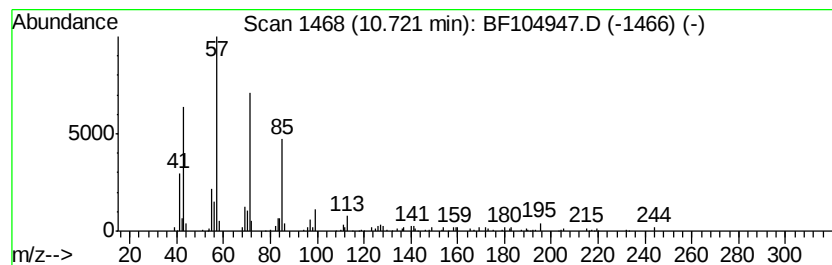
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Tridecane, 1-iodo- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.72	3.60 ng	157704	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
2	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
3	Eicosane	282	C20H42	000112-95-8	80
4	Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	72
5	Nonadecane	268	C19H40	000629-92-5	72



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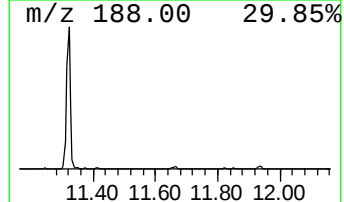
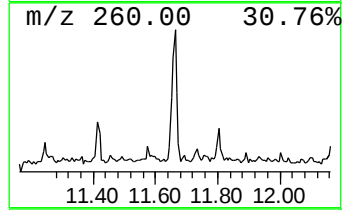
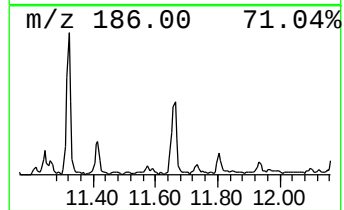
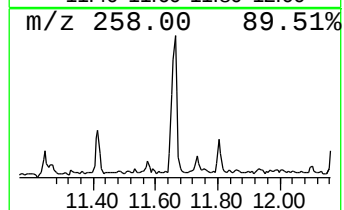
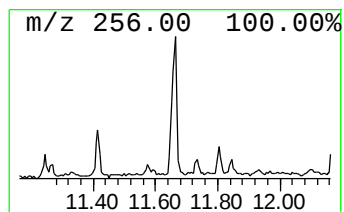
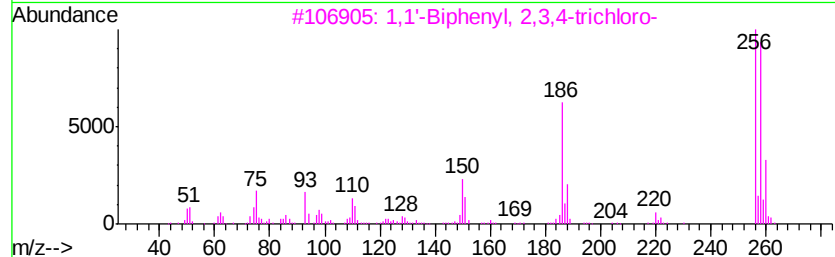
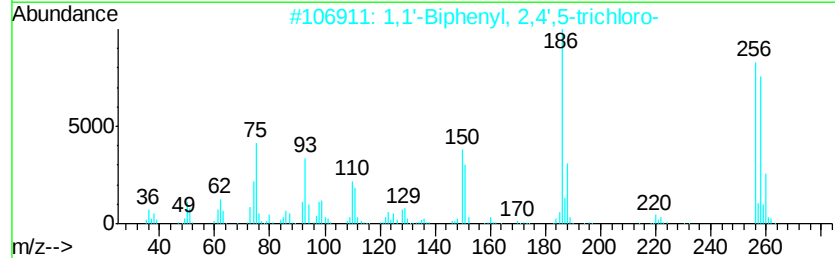
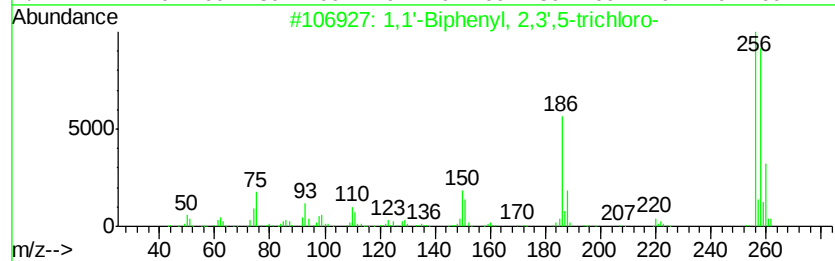
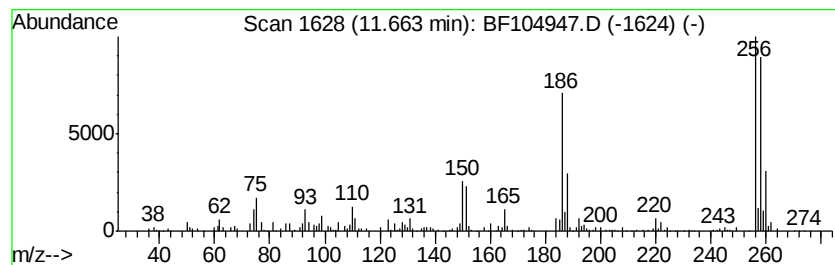
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2,3',5-trich... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	3.47 ng	151752	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	98
2			1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98
3			1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	98
4			1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98
5			1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	96



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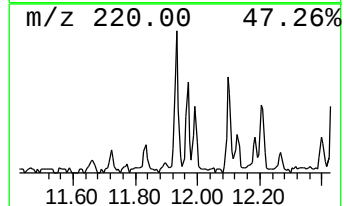
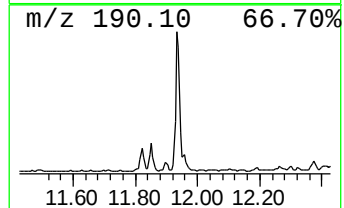
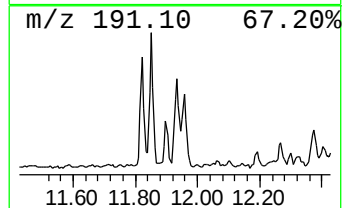
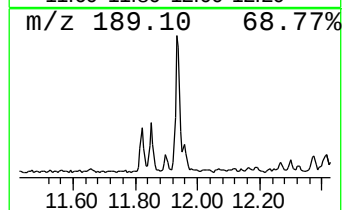
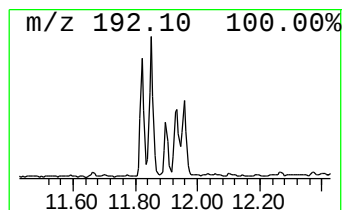
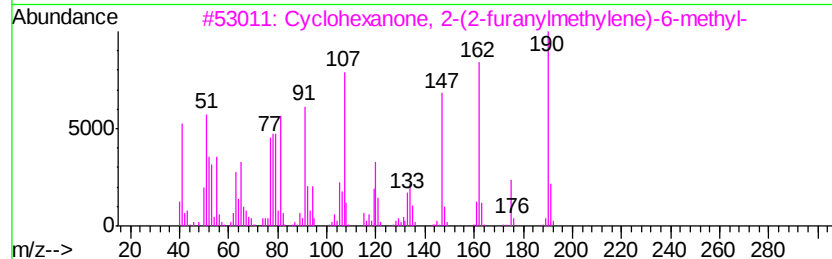
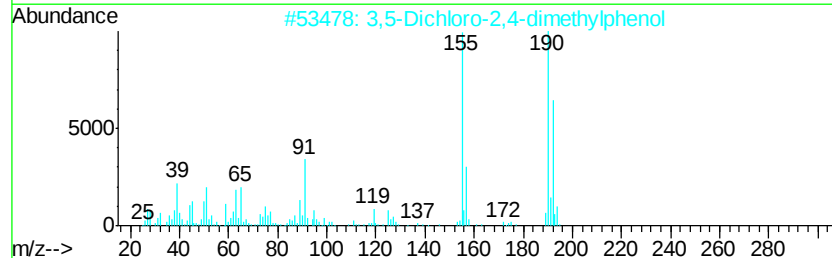
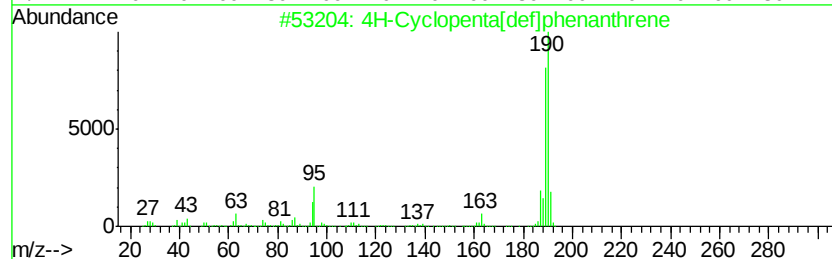
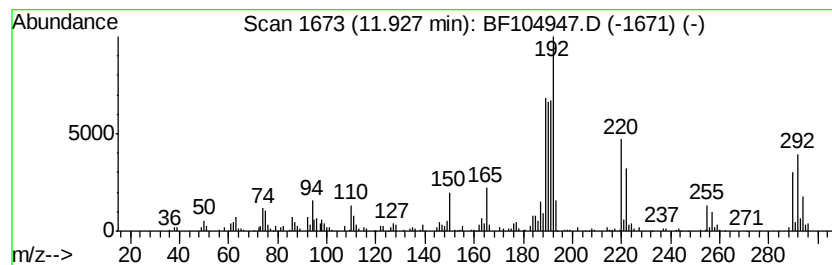
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Peak Number 5 unknown11.93 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.93	5.60 ng	245249	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
2	3,5-Dichloro-2,4-dimethylphenol	190	C8H8Cl2O	1000250-17-3	27
3	Cyclohexanone, 2-(2-furanylmethyl)-	190	C12H14O2	056053-07-7	22
4	Methyl 3,5-dichloro-4-hydroxybenzoate	220	C8H6Cl2O3	003337-59-5	22
5	3,5-Dichlorobenzoic acid	190	C7H4Cl2O2	000051-36-5	22



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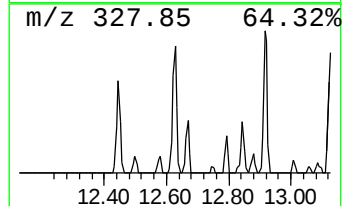
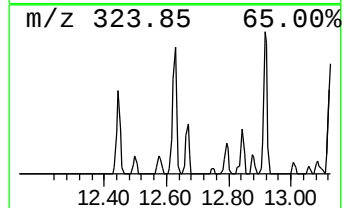
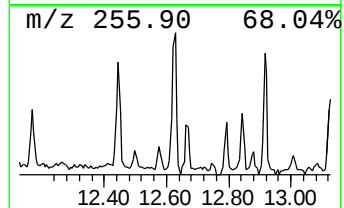
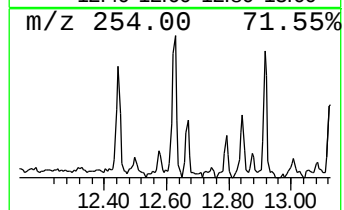
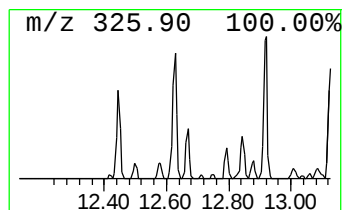
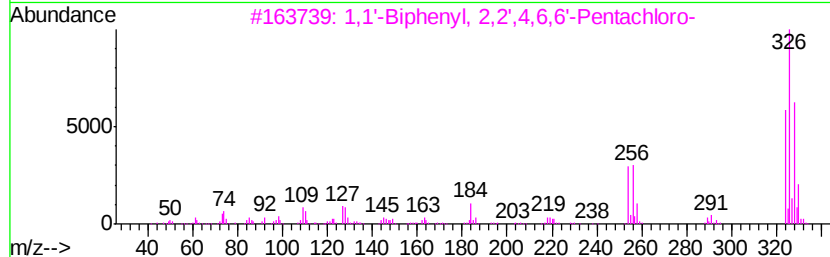
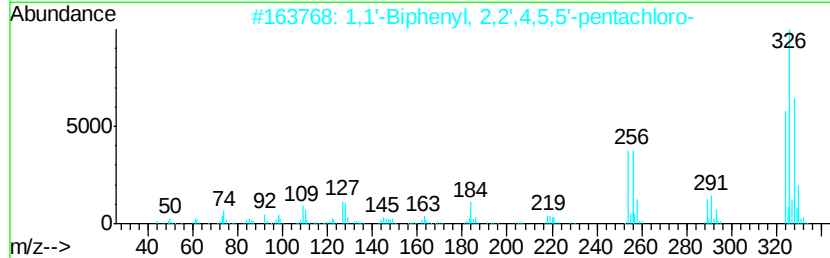
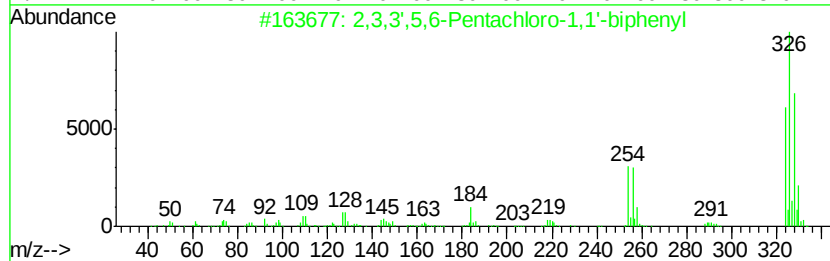
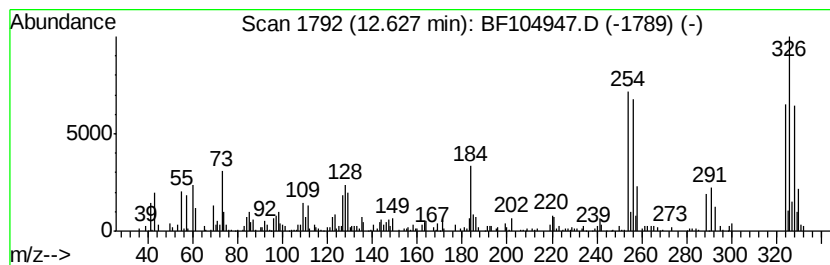
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Peak Number 6 2,3,3',5,6-Pentachloro-1,1'... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	3.62 ng	158442	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99		
2	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99		
3	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99		
4	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	99		
5	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	99		



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043018\
Data File : BF104947.D
Acq On : 30 Apr 2018 21:41
Operator : JU/SJ
Sample : J2609-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01A

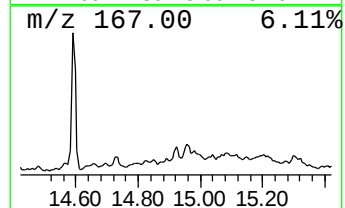
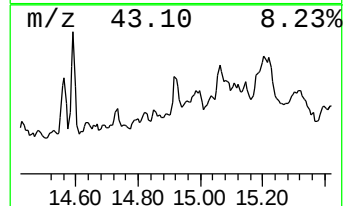
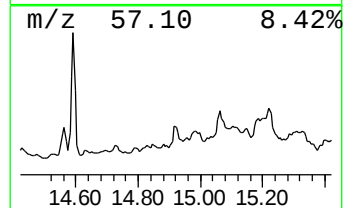
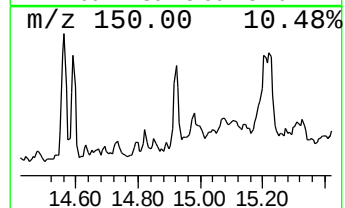
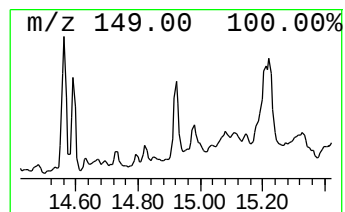
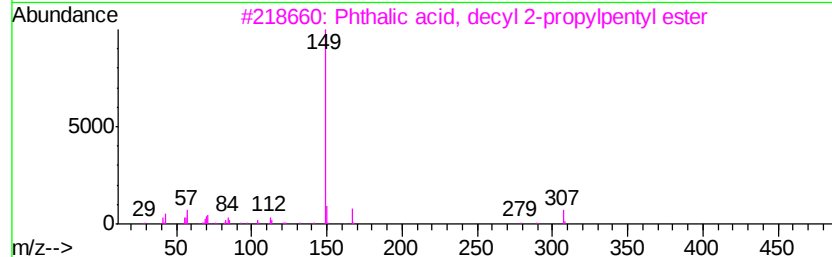
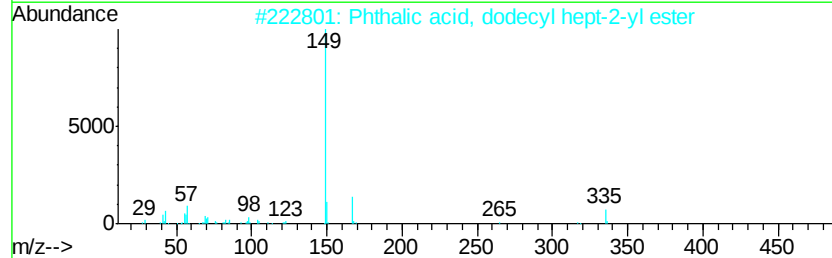
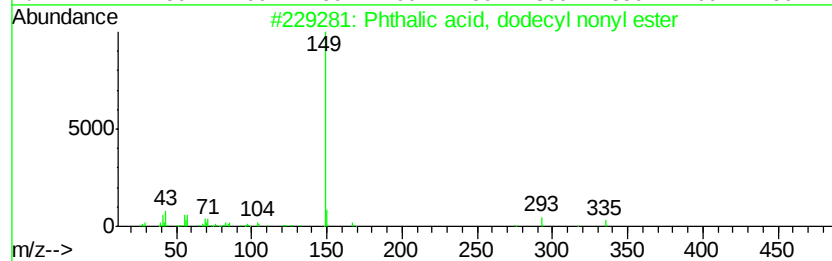
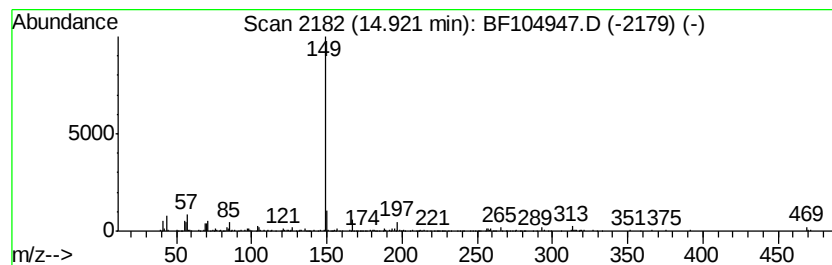
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phthalic acid, dodecyl nonyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	10.43 ng	242552	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	72
2		Phthalic acid, dodecyl hept-2-yl...	432	C27H44O4	1000356-97-9	72
3		Phthalic acid, decyl 2-propylpen...	418	C26H42O4	1000377-92-6	72
4		Phthalic acid, heptadecyl hept-4...	502	C32H54O4	1000356-79-8	72
5		Phthalic acid, 4-methylhept-3-yl...	404	C25H40O4	1000377-94-4	72



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01A

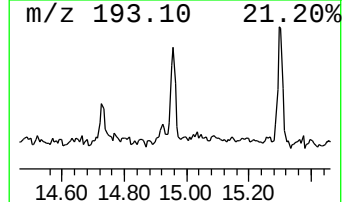
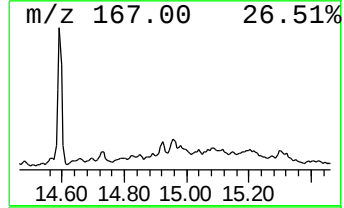
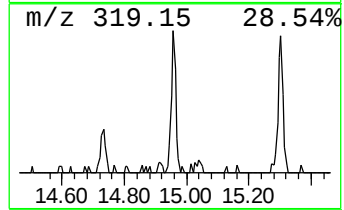
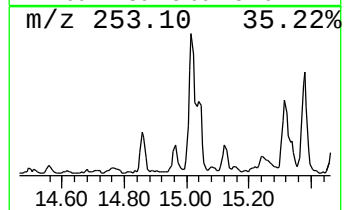
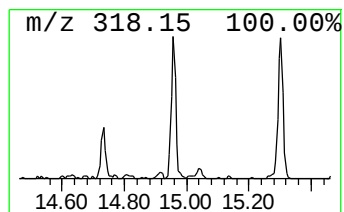
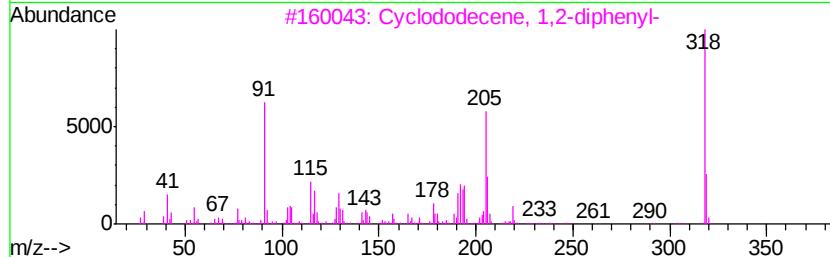
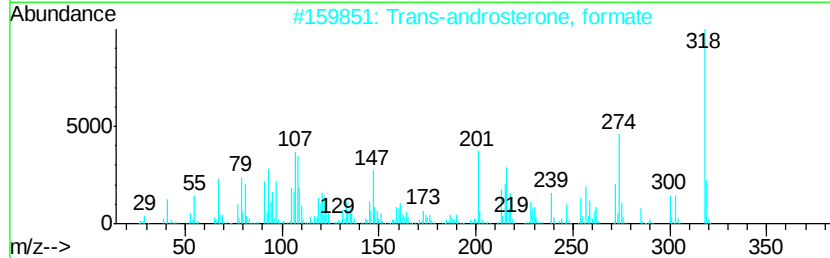
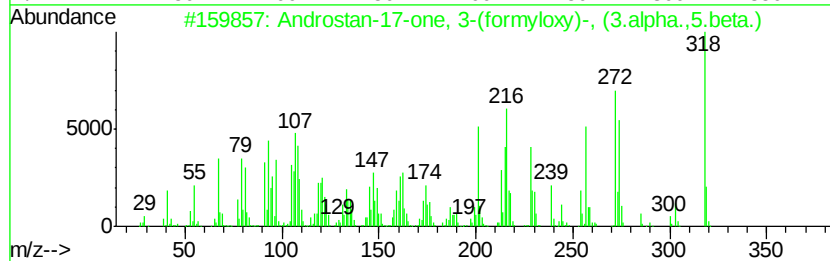
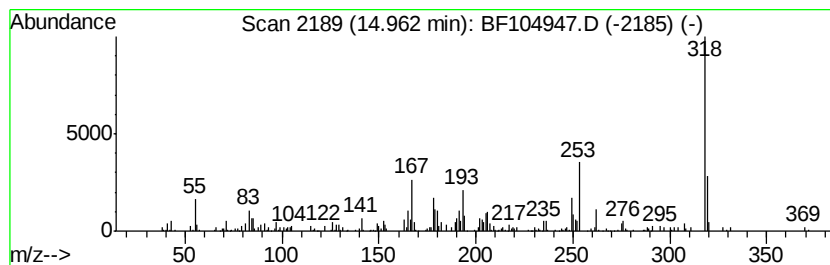
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.96 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	6.71 ng	156115	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androstan-17-one, 3-(formvloxy)-...	318	C20H30O3	004589-75-7	38
2		Trans-androsterone, formate	318	C20H30O3	1000368-37-9	38
3		Cyclododecene, 1,2-diphenyl-	318	C24H30	115977-56-5	38
4		2H-Chromene-3-carboxamide, 8-all...	318	C20H18N2O2	313497-20-0	30
5		Dibenzene, 2,2',5,5'-tetrafluoro...	318	C14H10F4O4	1000116-58-4	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043018\
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Acq On : 30 Apr 2018 21:41
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Sample : J2609-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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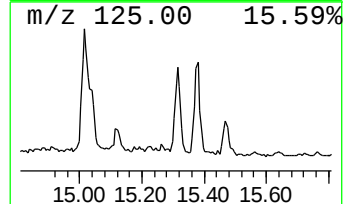
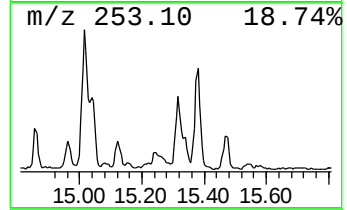
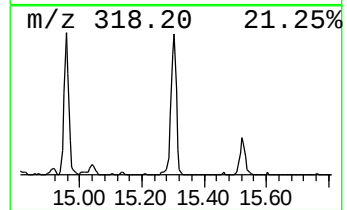
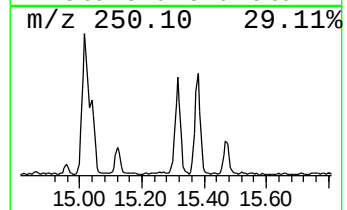
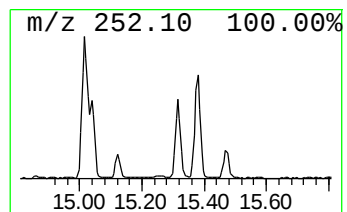
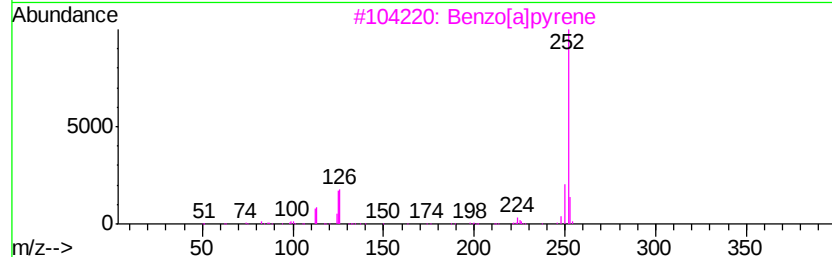
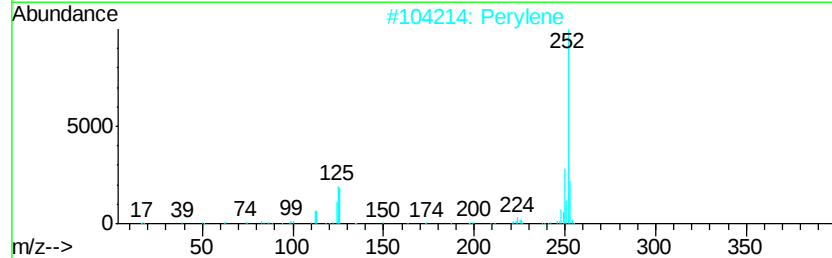
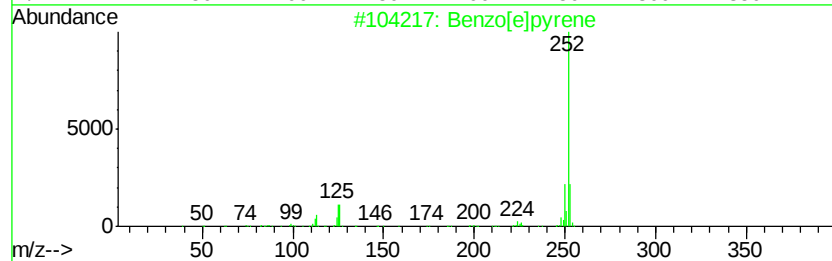
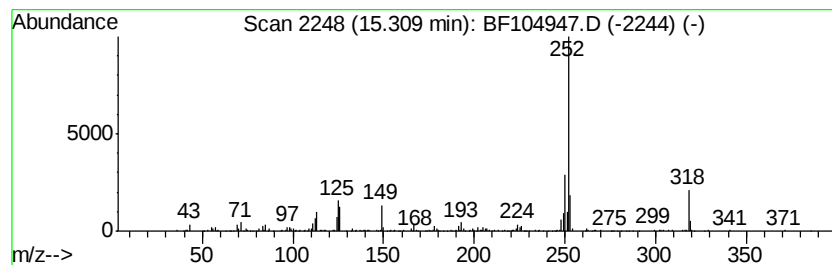
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.31	22.32 ng	518934	Perylene-d12	15.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	97
2		Perylene	252	C20H12	000198-55-0	96
3		Benzo[a]pyrene	252	C20H12	000050-32-8	87
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	78
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043018\
Data File : BF104947.D
Acq On : 30 Apr 2018 21:41
Operator : JU/SJ
Sample : J2609-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01A

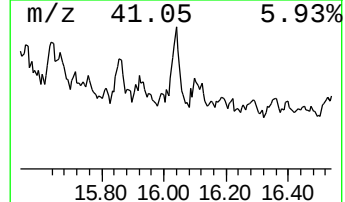
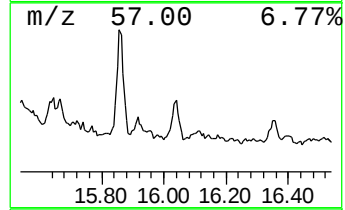
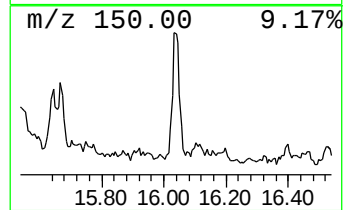
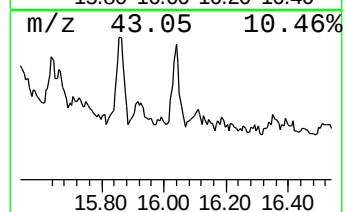
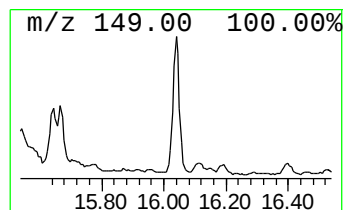
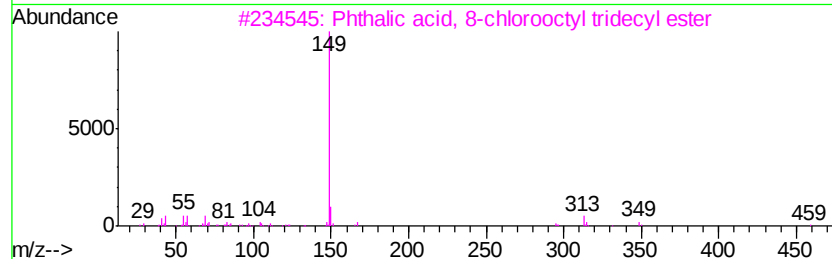
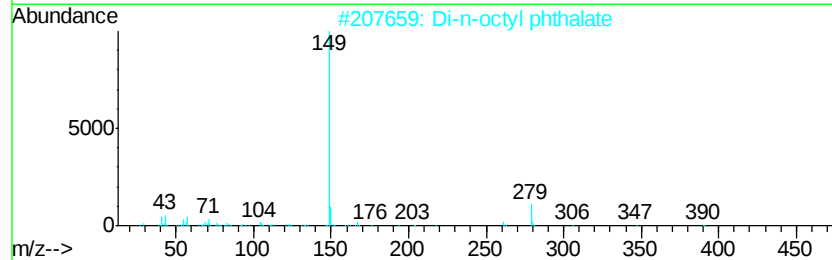
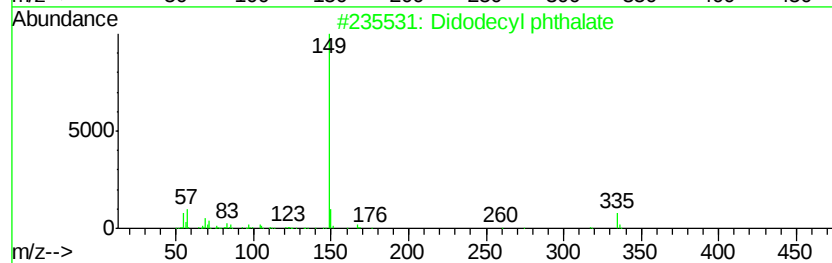
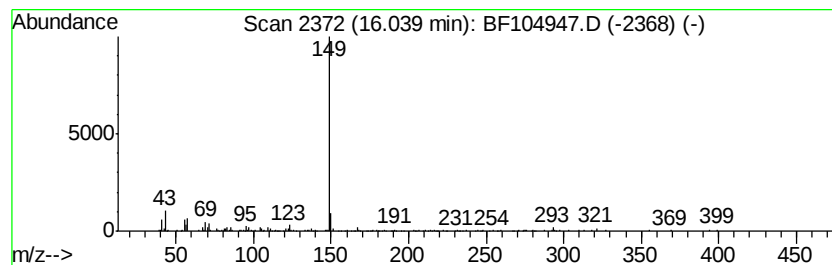
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Didodecyl phthalate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	10.34 ng	240324	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Didodecyl phthalate	502	C32H54O4	002432-90-8	80
2		Di-n-octyl phthalate	390	C24H38O4	000117-84-0	72
3		Phthalic acid, 8-chlorooctyl tri...	494	C29H47ClO4	1000356-86-8	72
4		Phthalic acid, cycloheptyl undec...	416	C26H40O4	1000322-83-5	72
5		Phthalic acid, hexadecyl hex-3-y...	474	C30H50O4	1000356-96-6	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043018\
Data File : BF104947.D
Acq On : 30 Apr 2018 21:41
Operator : JU/SJ
Sample : J2609-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

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ClientSampleId :
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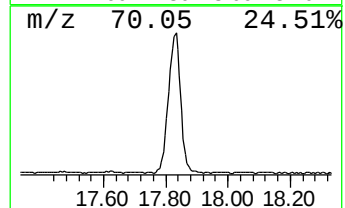
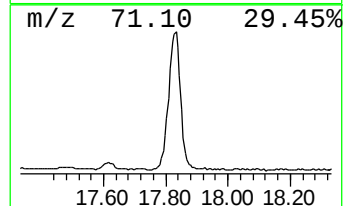
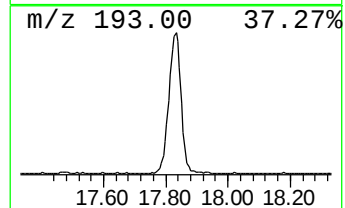
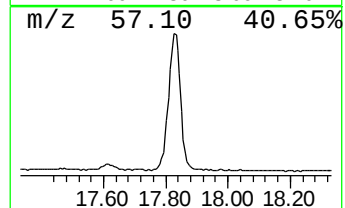
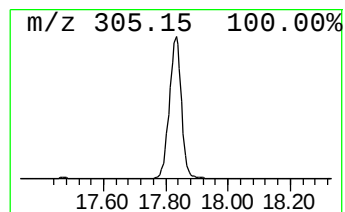
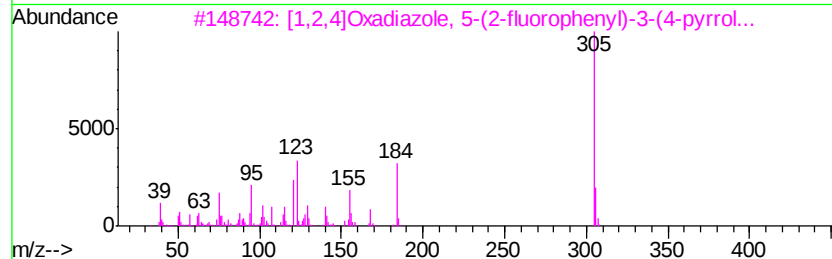
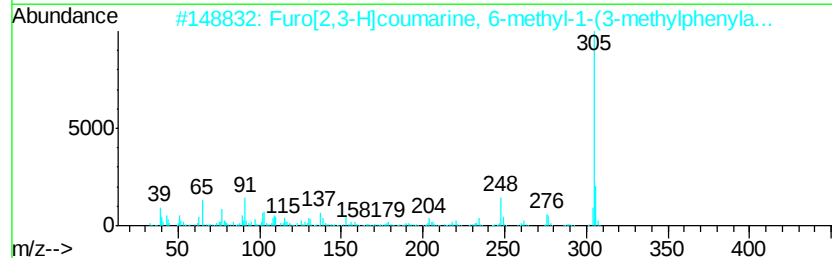
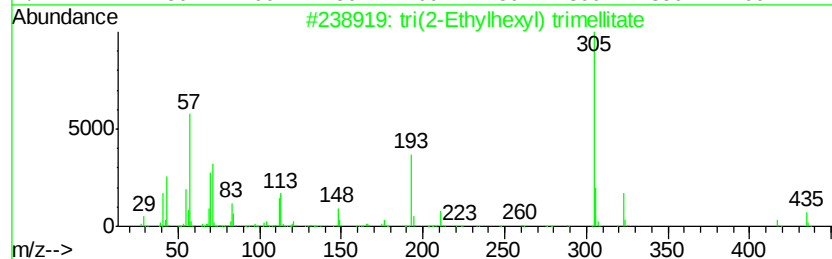
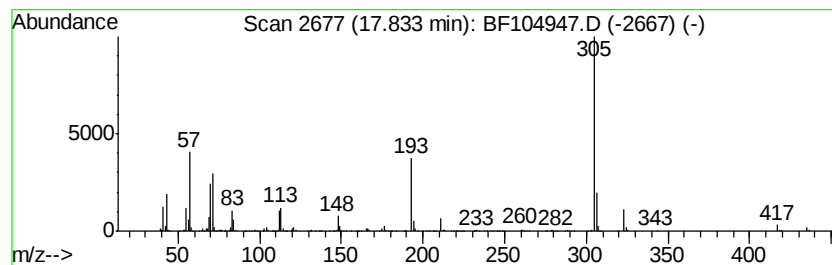
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 tri(2-Ethylhexyl) trimellitate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	95.30 ng	2215820	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	62
2		Furo[2,3-H]coumarine, 6-methyl-1...	305	C19H15N03	1000270-15-3	25
3		[1,2,4]Oxadiazole, 5-(2-fluoroph...	305	C18H12FN3O	1000316-03-9	9
4		Ethene, 1-(anthracen-9-yl)-2-(4-...	305	C23H15N	1000163-13-9	9
5		9H-Imidazo[1,2-a]benzimidazole, ...	305	C20H23N3	309740-57-6	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043018\
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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Tridecane, 1-iodo-	10.72	3.6	ng	157704	4	11.32	875772	20.0
1,1'-Biphenyl, 2,...	11.66	3.5	ng	151752	4	11.32	875772	20.0
unknown11.93	11.93	5.6	ng	245249	4	11.32	875772	20.0
2,3,3',5,6-Pentac...	12.63	3.6	ng	158442	4	11.32	875772	20.0
Phthalic acid, do...	14.92	10.4	ng	242552	6	15.44	465043	20.0
unknown14.96	14.96	6.7	ng	156115	6	15.44	465043	20.0
Benzo[e]pyrene	15.31	22.3	ng	518934	6	15.44	465043	20.0
Didodecyl phthalate	16.04	10.3	ng	240324	6	15.44	465043	20.0
tri(2-Ethylhexyl)...	17.83	95.3	ng	2215820	6	15.44	465043	20.0