

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF031618\
 Data File : BF103706.D
 Acq On : 16 Mar 2018 19:46
 Operator : JU/SJ
 Sample : J1929-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LS-PIT-2

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-BF031518.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	2.310	33	38	44	rVB	143327	184263	5.18%	0.536%
2	5.216	524	532	538	rBV	84223	102444	2.88%	0.298%
3	5.592	590	596	599	rBV	3548705	3495100	98.29%	10.173%
4	6.586	759	765	768	rBV	3182940	3555923	100.00%	10.350%
5	6.739	786	791	794	rBV	3705989	3540839	99.58%	10.306%
6	6.969	826	830	833	rBV	1320077	998944	28.09%	2.908%
7	7.128	852	857	860	rBV	3085892	2703987	76.04%	7.870%
8	7.528	920	925	929	rBV	2146802	2226985	62.63%	6.482%
9	8.251	1044	1048	1051	rBV	1702133	1346059	37.85%	3.918%
10	9.327	1225	1231	1234	rBV	3736485	3508823	98.68%	10.213%
11	9.392	1239	1242	1246	rBV	56690	57813	1.63%	0.168%
12	9.716	1293	1297	1299	rBV	316361	265651	7.47%	0.773%
13	9.927	1329	1333	1335	rVB	151544	126929	3.57%	0.369%
14	10.010	1343	1347	1350	rVB	1741537	1396803	39.28%	4.066%
15	10.245	1384	1387	1391	rBV2	36168	40866	1.15%	0.119%
16	10.427	1415	1418	1421	rVB	194218	156925	4.41%	0.457%
17	10.645	1450	1455	1458	rBV	59998	71948	2.02%	0.209%
18	10.798	1477	1481	1484	rBV	2140782	1965383	55.27%	5.720%
19	10.898	1495	1498	1505	rVB	145860	159690	4.49%	0.465%
20	11.057	1522	1525	1528	rVB	57656	49926	1.40%	0.145%
21	11.257	1556	1559	1565	rVB3	44068	46363	1.30%	0.135%
22	11.351	1572	1575	1577	rBV	73491	63396	1.78%	0.185%
23	11.380	1577	1580	1585	rVB3	39987	49231	1.38%	0.143%
24	11.498	1597	1600	1603	rBV	1331798	1222244	34.37%	3.557%
25	11.521	1603	1604	1610	rVB2	89352	81119	2.28%	0.236%
26	12.010	1683	1687	1692	rBV2	160975	163684	4.60%	0.476%
27	12.716	1802	1807	1810	rBV	146034	149710	4.21%	0.436%
28	12.774	1812	1817	1819	rVV2	131241	174100	4.90%	0.507%
29	12.798	1819	1821	1827	rVB4	43281	54527	1.53%	0.159%
30	12.898	1835	1838	1840	rVB2	48228	41544	1.17%	0.121%
31	12.945	1840	1846	1849	rBV2	142291	161903	4.55%	0.471%
32	13.021	1856	1859	1863	rVB	446479	406982	11.45%	1.185%
33	13.086	1866	1870	1874	rVB	2688552	2607456	73.33%	7.589%
34	13.280	1900	1903	1906	rBV4	39163	36659	1.03%	0.107%

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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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.363	1911	1917	1920	rBV	857008	732859	20.61%	2.133%
36	13.610	1956	1959	1963	rBV2	35691	61026	1.72%	0.178%
37	13.663	1965	1968	1971	rVV	69765	68655	1.93%	0.200%
38	13.715	1974	1977	1981	rVV	85417	77427	2.18%	0.225%
39	13.968	2017	2020	2025	rBV	380449	380095	10.69%	1.106%
40	14.151	2047	2051	2054	rBV	946934	906063	25.48%	2.637%
41	15.686	2307	2312	2327	rVB	637440	916710	25.78%	2.668%

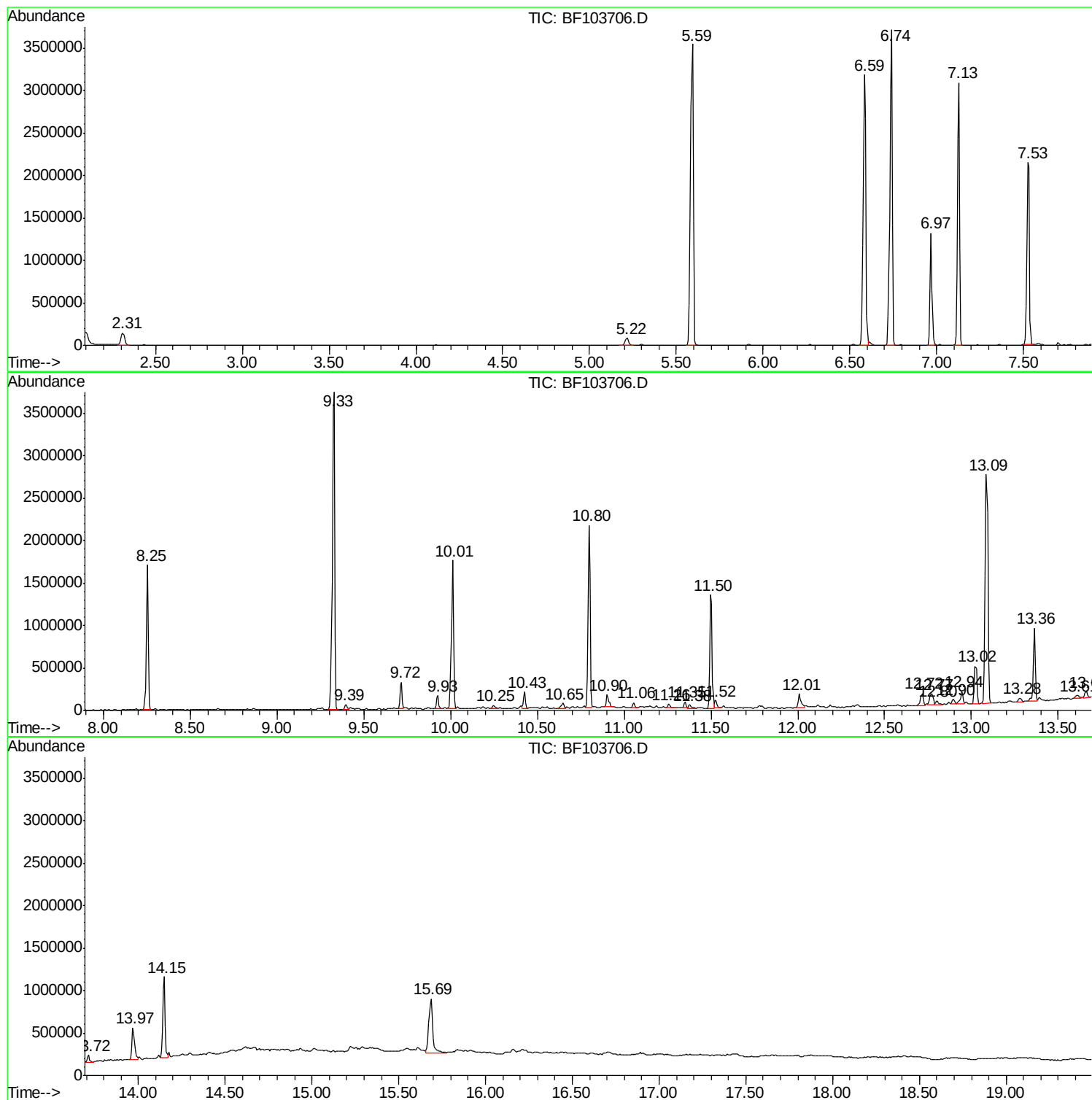
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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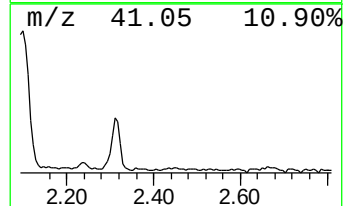
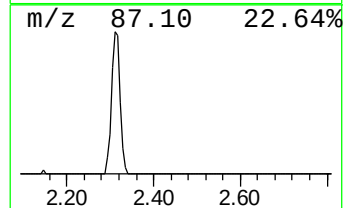
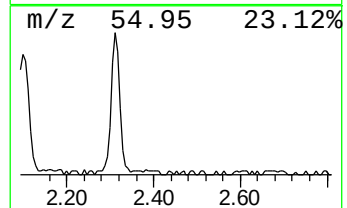
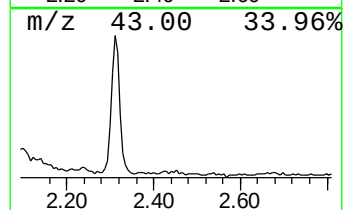
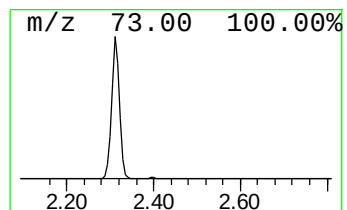
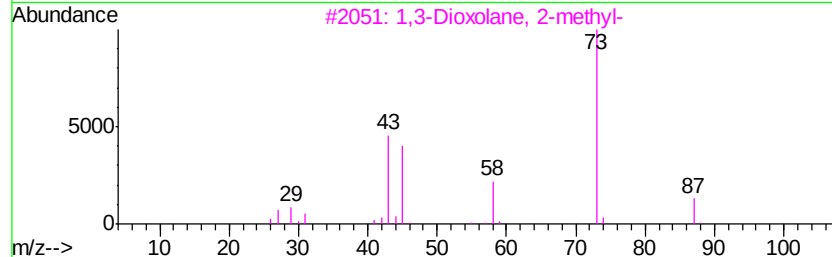
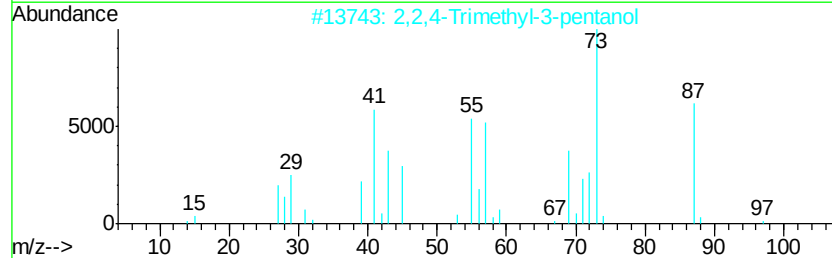
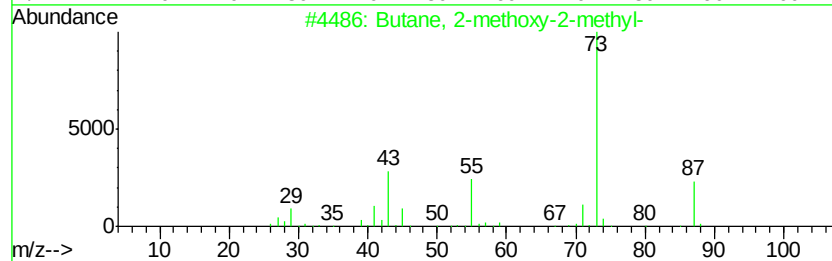
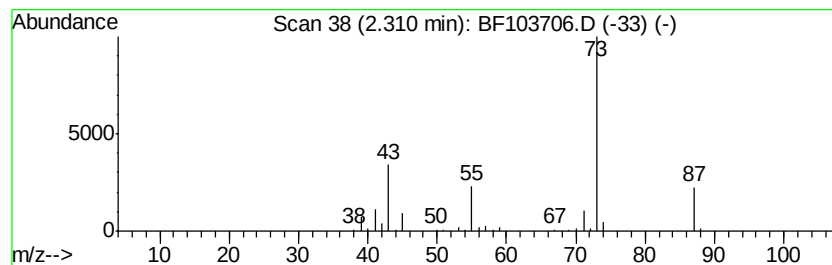
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.31	3.69 ng	184263	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17



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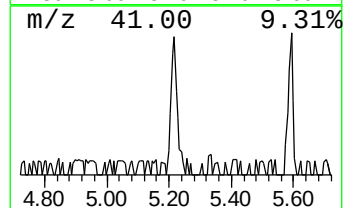
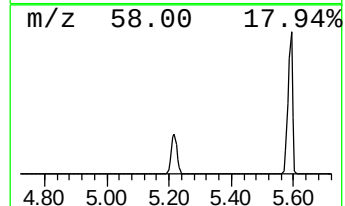
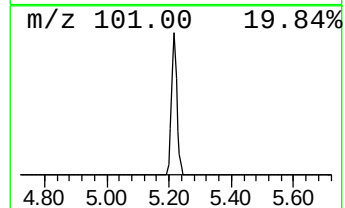
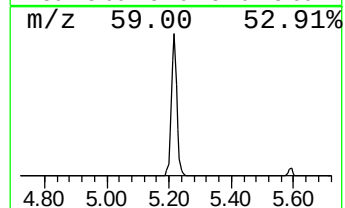
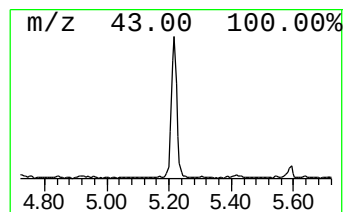
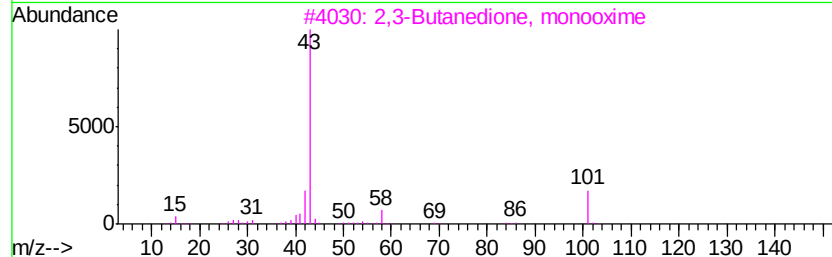
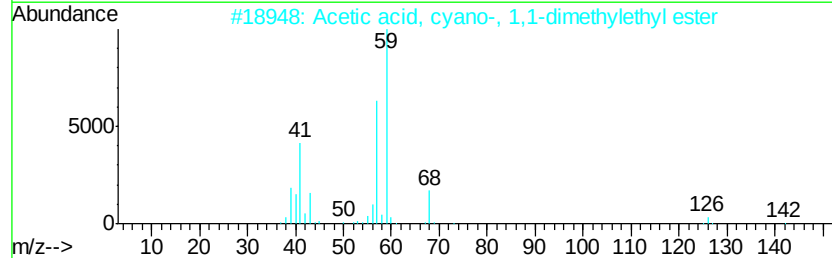
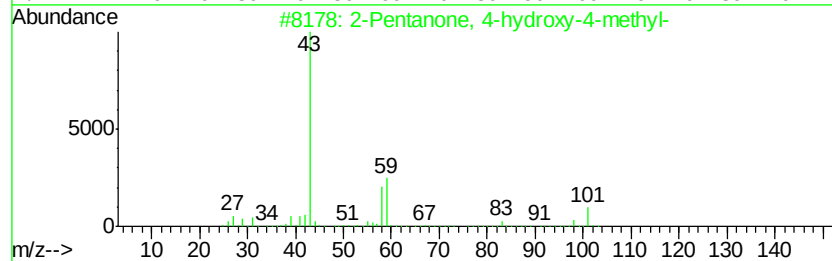
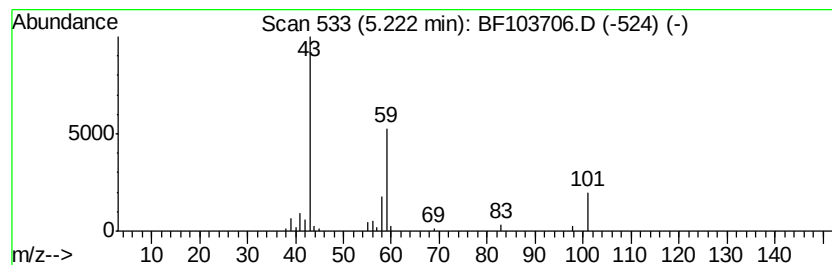
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	2.05 ng	102444	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	066241-43-8	9



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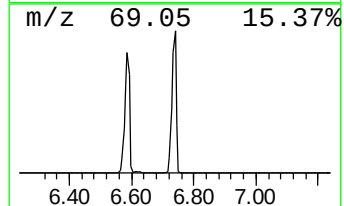
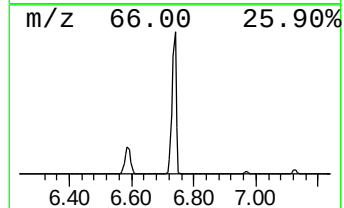
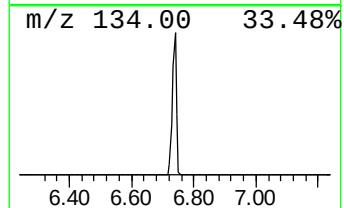
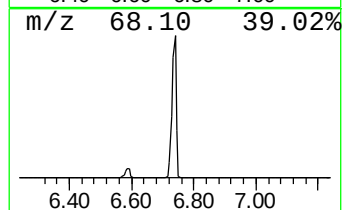
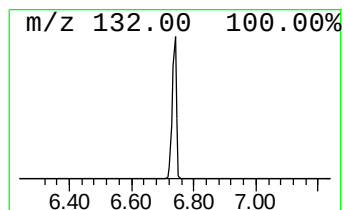
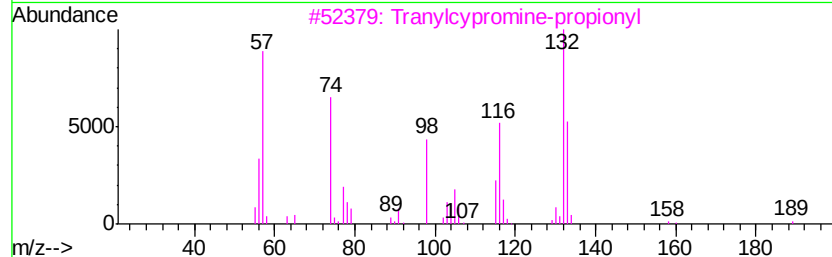
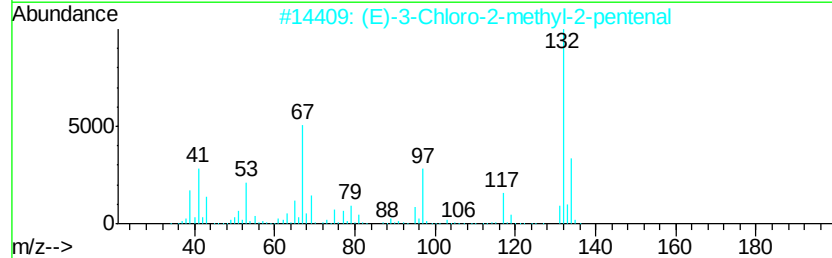
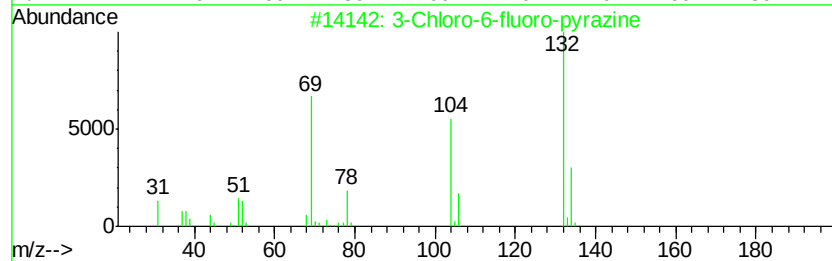
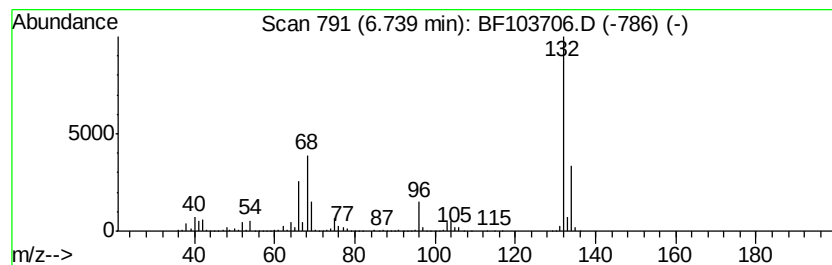
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	70.89 ng	3540840	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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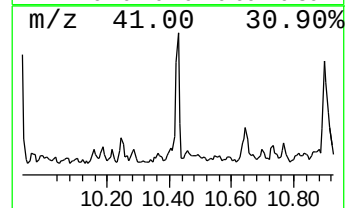
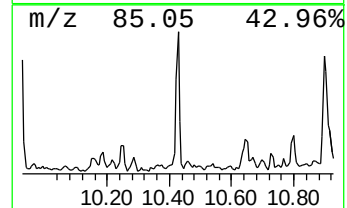
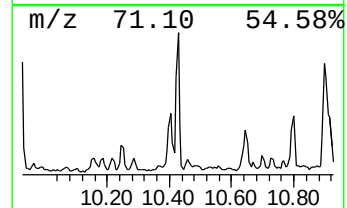
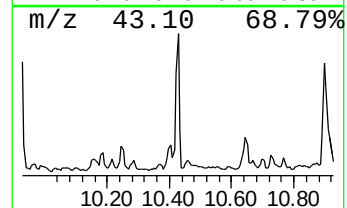
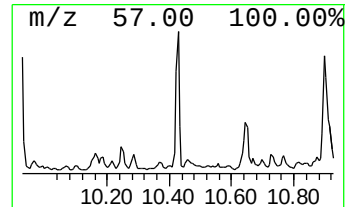
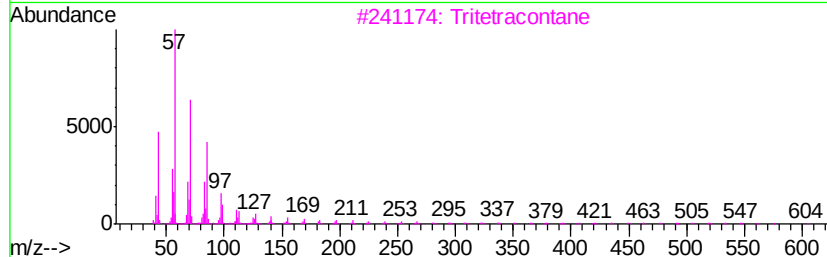
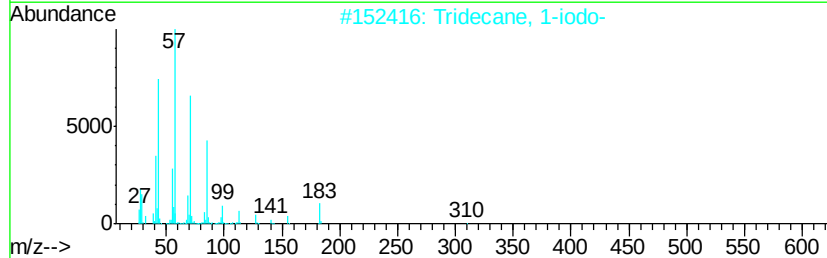
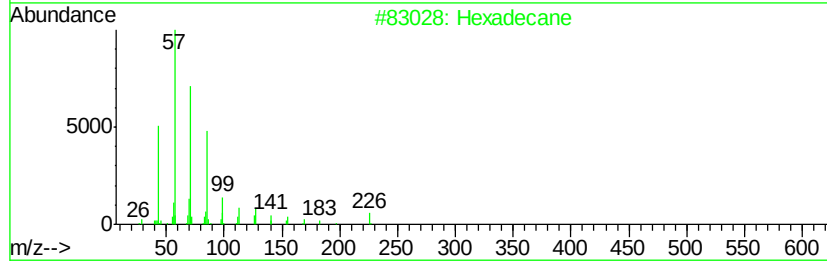
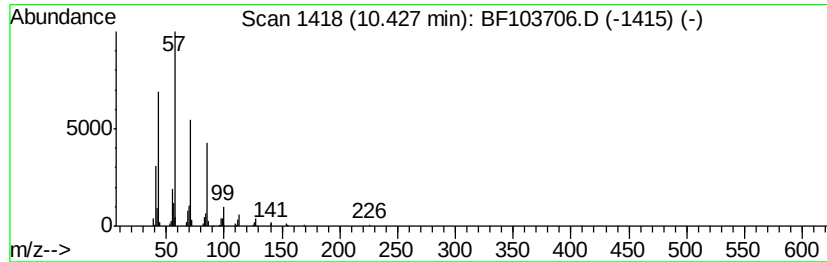
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.43	2.25 ng	156925	Acenaphthene-d10	10.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3	Tritetracontane	605	C43H88	007098-21-7	90
4	Heptadecane	240	C17H36	000629-78-7	90
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90



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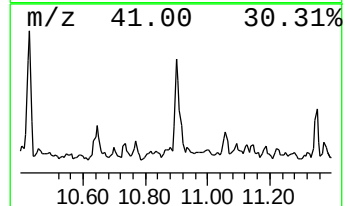
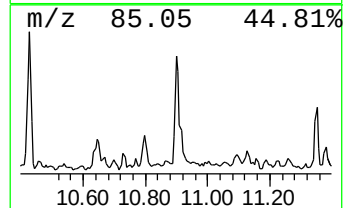
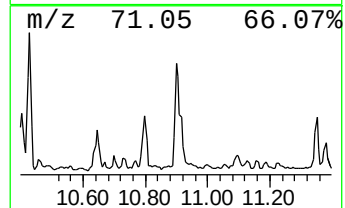
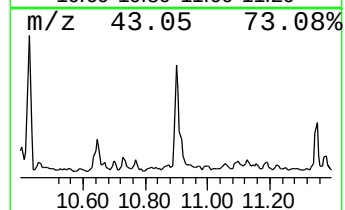
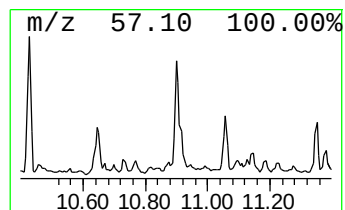
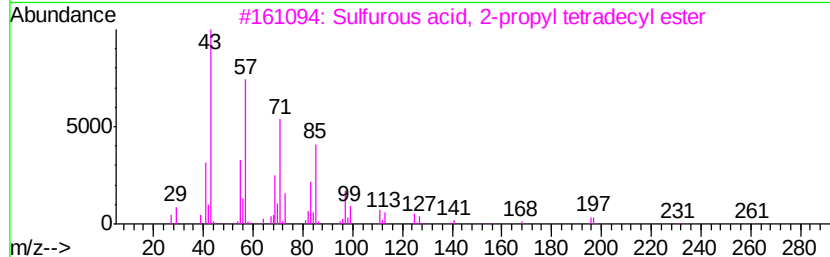
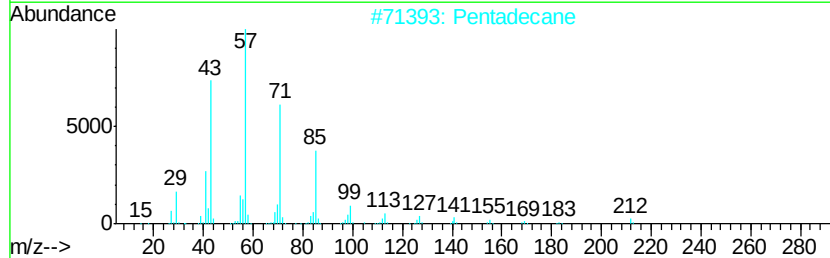
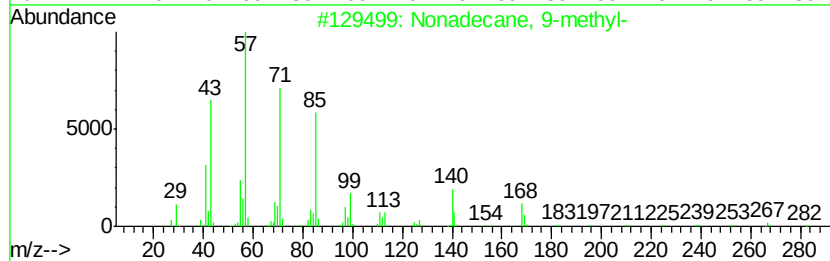
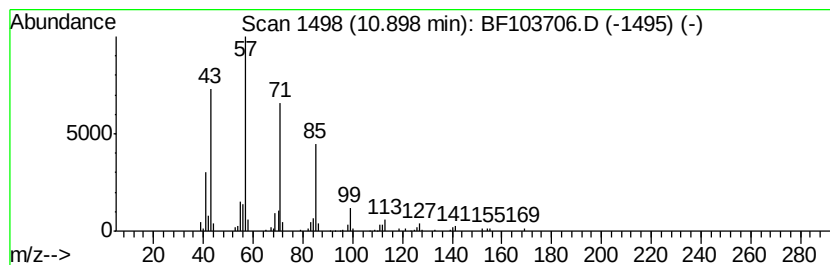
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Nonadecane, 9-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	2.61 ng	159690	Phenanthrene-d10	11.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	90
2	Pentadecane	212	C15H32	000629-62-9	90
3	Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	78
4	Hexadecane	226	C16H34	000544-76-3	78
5	Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	78



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 Acq On : 16 Mar 2018 19:46
 Operator : JU/SJ
 Sample : J1929-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LS-PIT-2

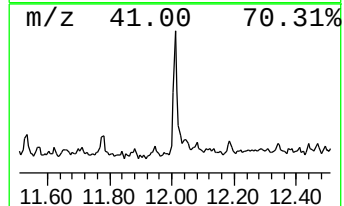
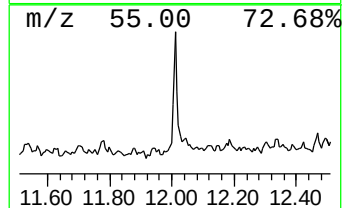
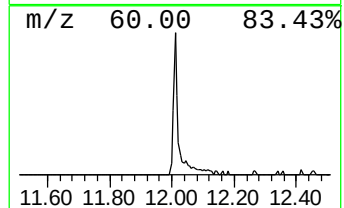
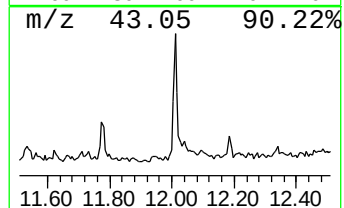
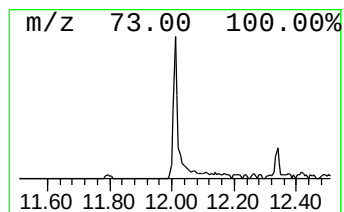
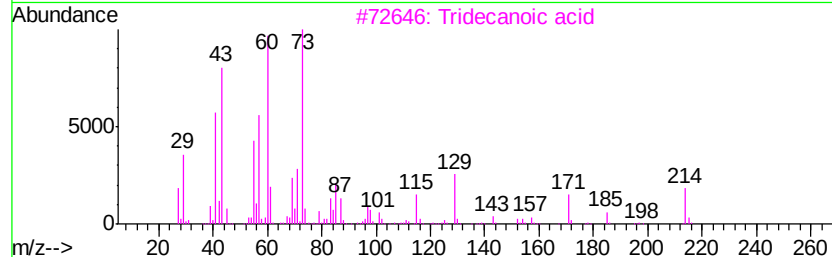
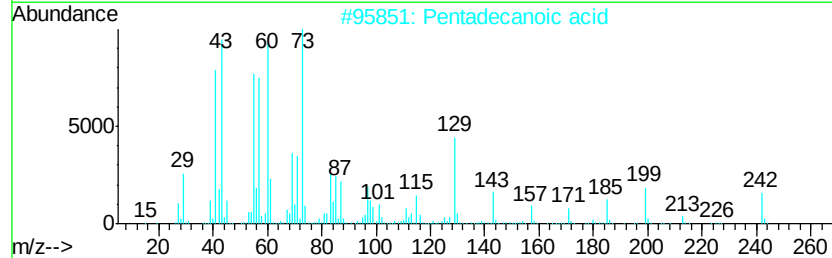
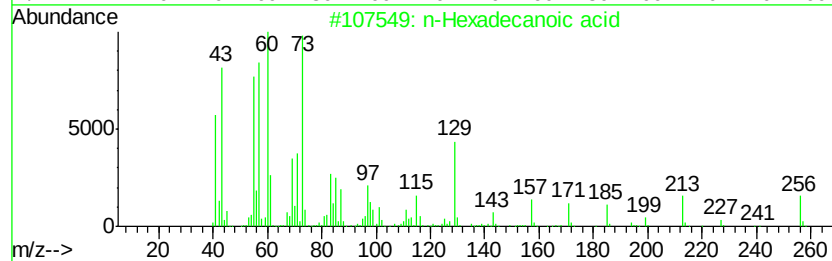
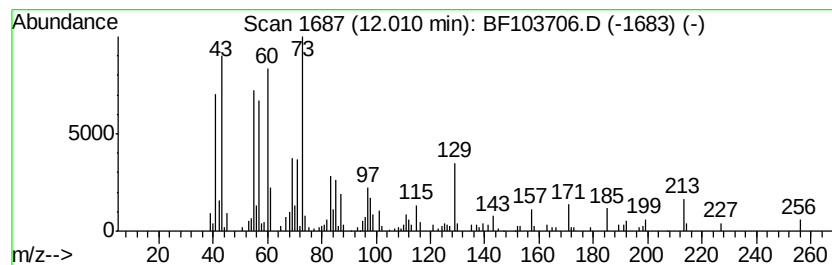
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	2.68 ng	163684	Phenanthrene-d10	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		Dodecanoic acid	200	C12H24O2	000143-07-7	53



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF031618\
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Operator : JU/SJ
Sample : J1929-05
Misc :
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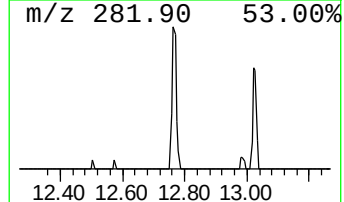
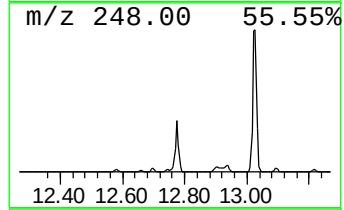
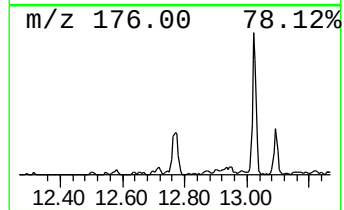
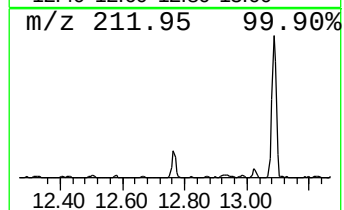
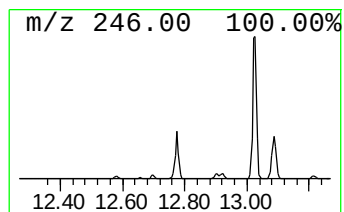
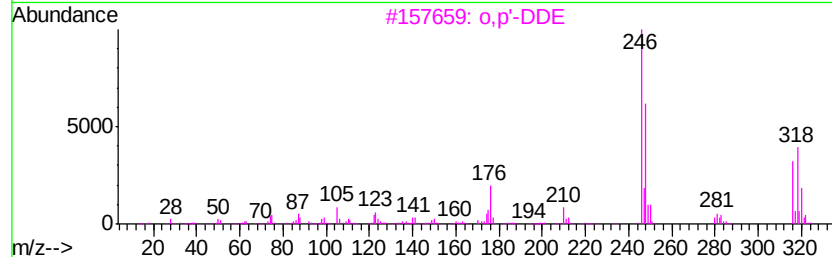
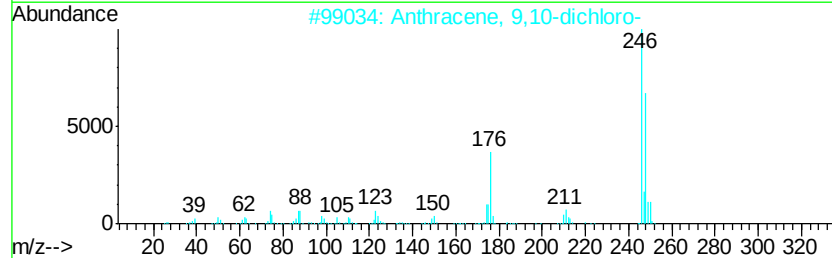
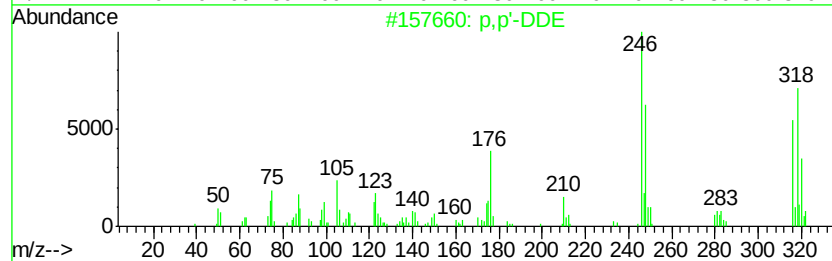
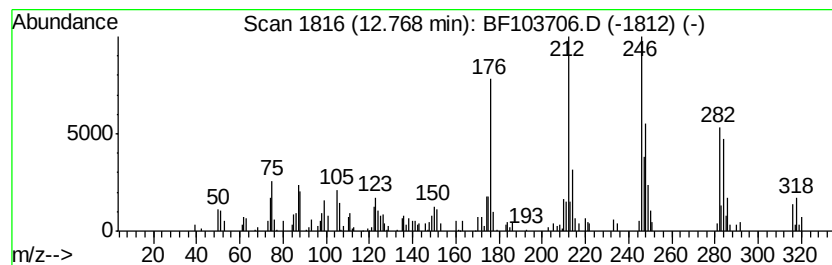
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 p,p'-DDE Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.77	2.85 ng	174100	Phenanthrene-d10	11.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDE	316	C14H8Cl4	000072-55-9	99
2	Anthracene, 9,10-dichloro-	246	C14H8Cl2	000605-48-1	98
3	o,p'-DDE	316	C14H8Cl4	003424-82-6	98
4	Anthracene, 1,8-dichloro-	246	C14H8Cl2	014381-66-9	93
5	9H-Fluorene, 9-(dichloromethylene)-	246	C14H8Cl2	000835-17-6	91



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF031618\
Data File : BF103706.D
Acq On : 16 Mar 2018 19:46
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Sample : J1929-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

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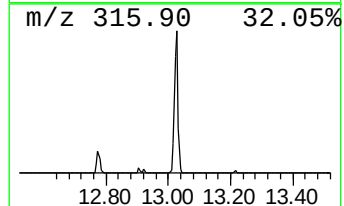
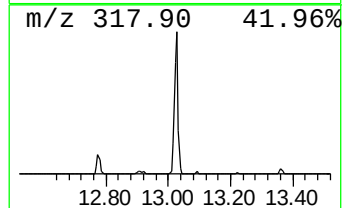
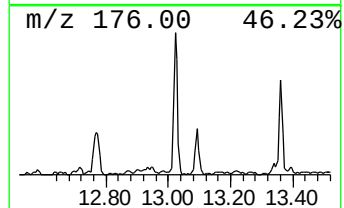
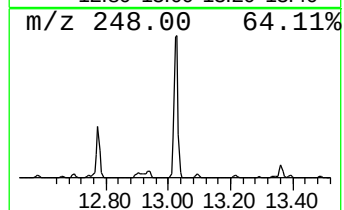
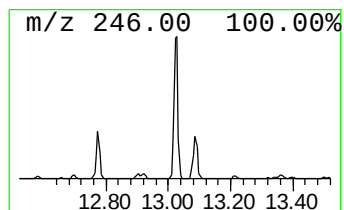
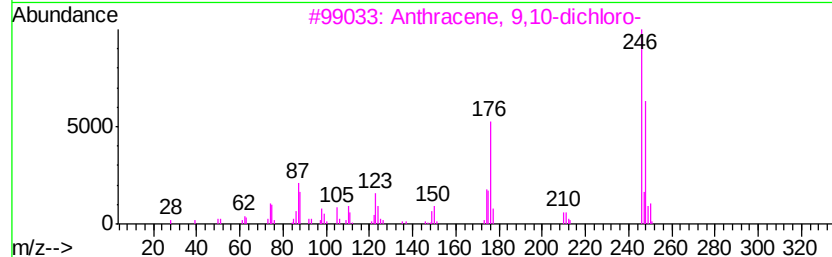
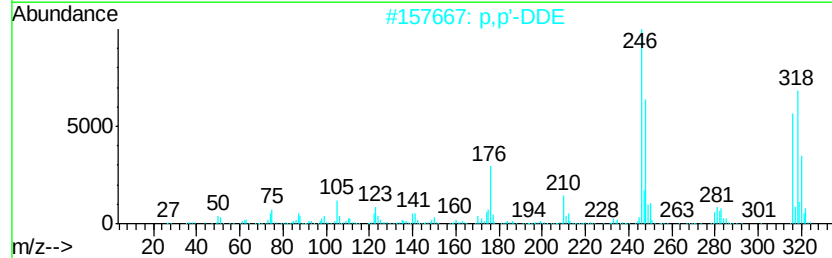
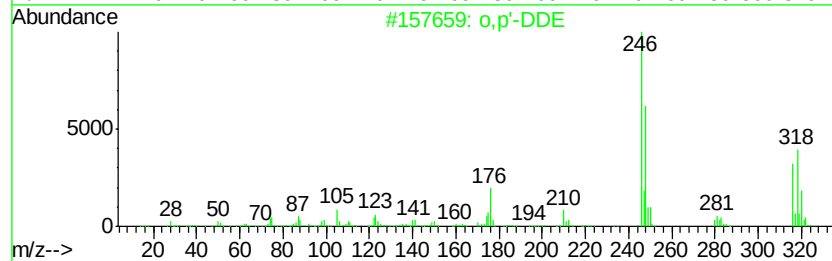
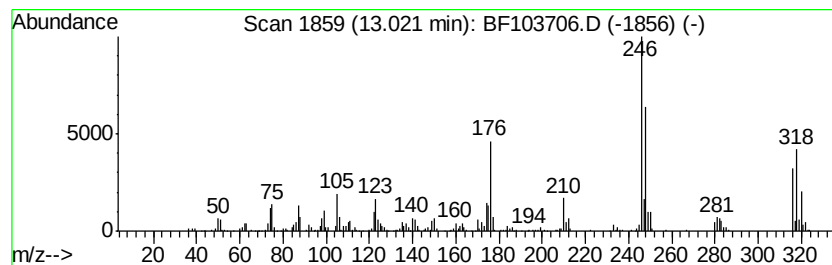
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 o,p'-DDE Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	8.98 ng	406982	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
2		p,p'-DDE	316	C14H8Cl4	000072-55-9	99
3		Anthracene, 9,10-dichloro-	246	C14H8Cl2	000605-48-1	98
4		Anthracene, 1,8-dichloro-	246	C14H8Cl2	014381-66-9	92
5		Benzene, 1,1'-(1,2-dichloro-1,2-...	316	C14H8Cl4	088970-69-8	91



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF031618\
Data File : BF103706.D
Acq On : 16 Mar 2018 19:46
Operator : JU/SJ
Sample : J1929-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

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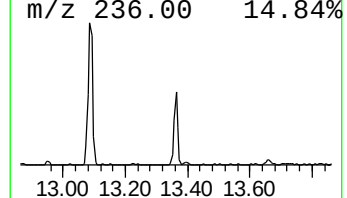
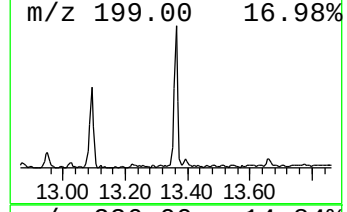
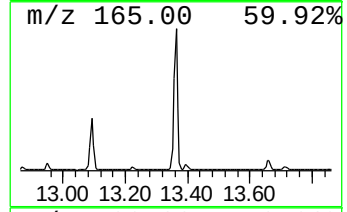
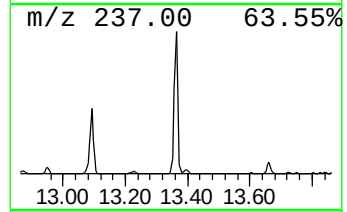
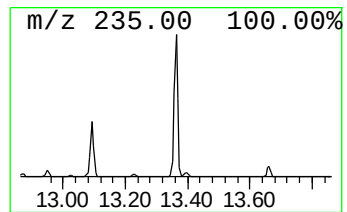
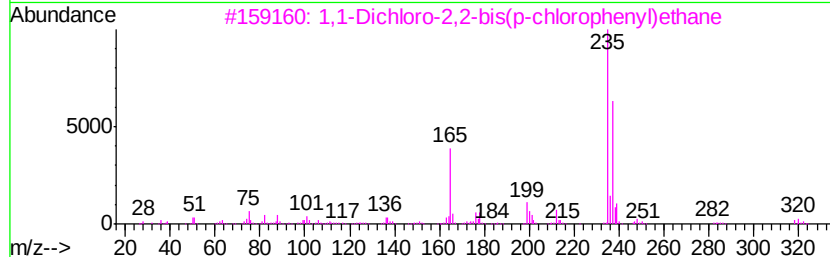
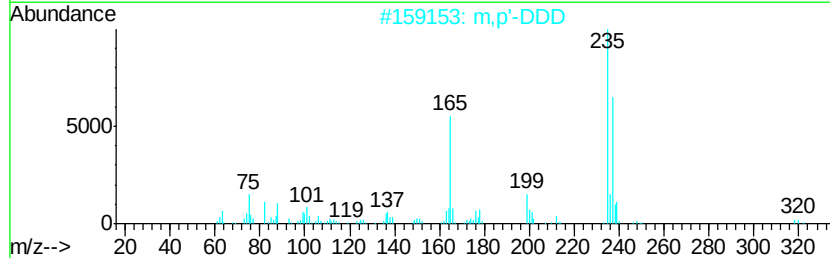
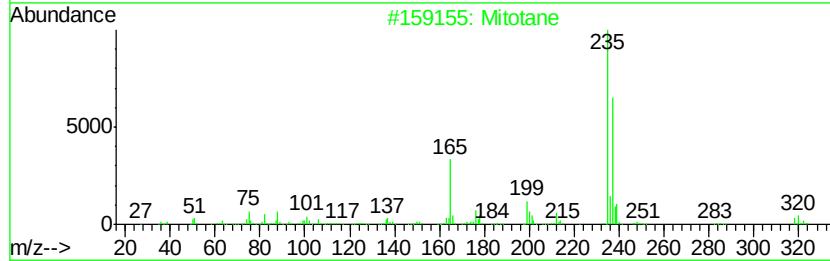
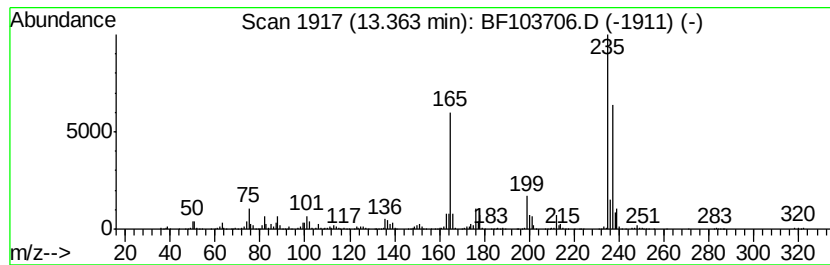
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Mitotane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	16.18 ng	732859	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mitotane	318	C14H10Cl4	000053-19-0	99
2		m,p'-DDD	318	C14H10Cl4	004329-12-8	99
3		1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	96
4		2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	86
5		3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	72



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Sample : J1929-05
Misc :
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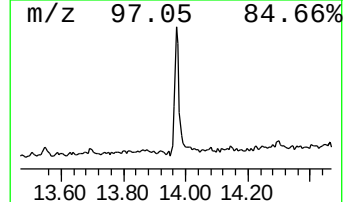
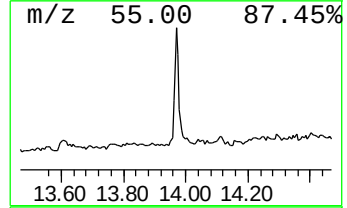
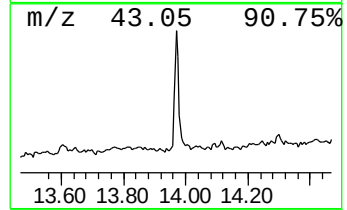
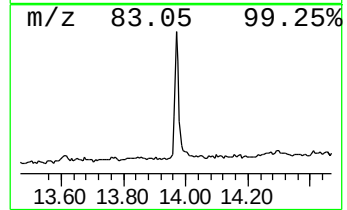
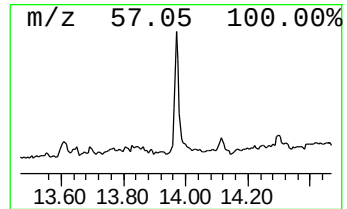
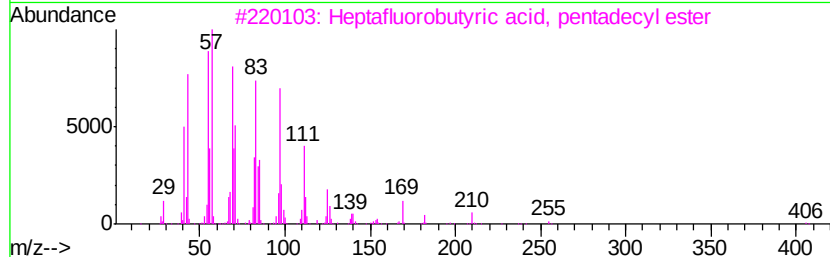
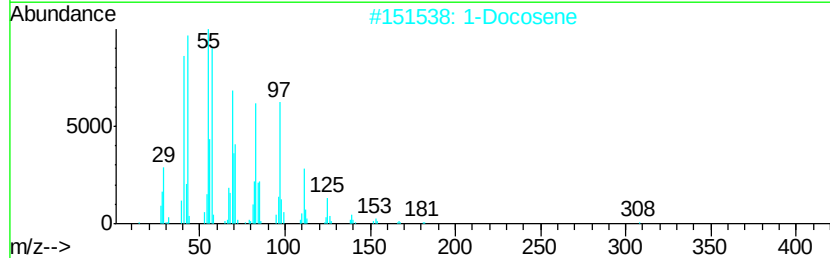
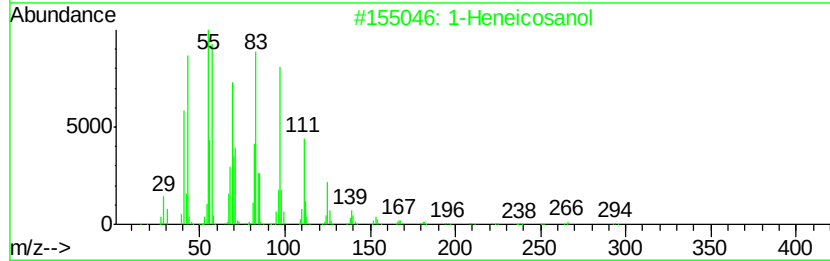
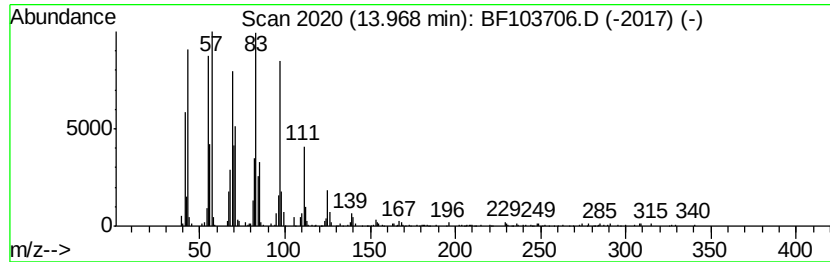
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.97	8.39 ng	380095	Chrysene-d12	14.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1-Docosene	308	C22H44	001599-67-3	94
3	Heptafluorobutvric acid, pentade...	424	C19H31F7O2	959261-23-5	93
4	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF031618\
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Acq On : 16 Mar 2018 19:46
Operator : JU/SJ
Sample : J1929-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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2-Pentanone, 4-hy...	5.22	2.0	ng	102444	1	6.97	998944	20.0
unknown6.74	6.74	70.9	ng	3540840	1	6.97	998944	20.0
Hexadecane	10.43	2.3	ng	156925	3	10.01	1396800	20.0
Nonadecane, 9-met...	10.90	2.6	ng	159690	4	11.50	1222240	20.0
n-Hexadecanoic acid	12.01	2.7	ng	163684	4	11.50	1222240	20.0
p,p'-DDE	12.77	2.9	ng	174100	4	11.50	1222240	20.0
o,p'-DDE	13.02	9.0	ng	406982	5	14.15	906063	20.0
Mitotane	13.36	16.2	ng	732859	5	14.15	906063	20.0
1-Heneicosanol	13.97	8.4	ng	380095	5	14.15	906063	20.0