

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF112018\
 Data File : BF110904.D
 Acq On : 20 Nov 2018 19:45
 Operator : JU/SJ
 Sample : J5996-25 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOC1-06-A

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.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	2.187	13	17	30	rVB4	15686	29015	6.74%	0.879%
2	5.163	519	523	526	rBV	9833	12485	2.90%	0.378%
3	5.575	589	593	604	rBV	24363	55210	12.83%	1.673%
4	6.575	760	763	784	rBV	58380	122826	28.54%	3.723%
5	6.716	784	787	798	rVB	100567	112807	26.22%	3.419%
6	6.939	822	825	843	rBV	323898	334477	77.73%	10.137%
7	7.098	849	852	863	rBV	103231	113017	26.27%	3.425%
8	7.522	920	924	940	rBV	37931	76479	17.77%	2.318%
9	8.222	1040	1043	1064	rBV	318253	430294	100.00%	13.041%
10	9.222	1211	1213	1215	rVB	6713	4702	1.09%	0.143%
11	9.298	1223	1226	1234	rBV	117112	132161	30.71%	4.005%
12	9.363	1234	1237	1241	rVB2	9369	10662	2.48%	0.323%
13	9.680	1287	1291	1295	rBV2	19925	21332	4.96%	0.647%
14	9.733	1297	1300	1304	rVB3	3775	5187	1.21%	0.157%
15	9.833	1314	1317	1319	rBV	6191	5168	1.20%	0.157%
16	9.880	1323	1325	1330	rVV3	8160	9482	2.20%	0.287%
17	9.986	1339	1343	1355	rBV	371167	406587	94.49%	12.323%
18	10.392	1408	1412	1415	rBV	20317	18490	4.30%	0.560%
19	10.610	1445	1449	1452	rBV	11018	10915	2.54%	0.331%
20	10.792	1477	1480	1485	rVB4	7601	11329	2.63%	0.343%
21	10.875	1489	1494	1497	rVV	21017	22608	5.25%	0.685%
22	11.339	1570	1573	1576	rVB2	10456	9037	2.10%	0.274%
23	11.480	1593	1597	1609	rBV	369688	390399	90.73%	11.832%
24	11.733	1638	1640	1646	rVB3	5525	8083	1.88%	0.245%
25	11.980	1679	1682	1686	rBV3	5828	9610	2.23%	0.291%
26	12.104	1699	1703	1708	rBV6	4792	8531	1.98%	0.259%
27	12.533	1774	1776	1782	rBV4	13989	17820	4.14%	0.540%
28	12.704	1802	1805	1811	rBV	29017	33674	7.83%	1.021%
29	12.910	1838	1840	1842	rBV2	7844	8254	1.92%	0.250%
30	12.939	1842	1845	1851	rVV2	27866	50409	11.72%	1.528%
31	13.057	1862	1865	1871	rBV	155447	158669	36.87%	4.809%
32	13.357	1914	1916	1918	rBV3	12805	10618	2.47%	0.322%
33	14.133	2045	2048	2055	rBV	345298	370913	86.20%	11.241%
34	14.868	2171	2173	2176	rVB	43164	41144	9.56%	1.247%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.668	2305	2309	2318	rVB	166377	237135	55.11%	7.187%
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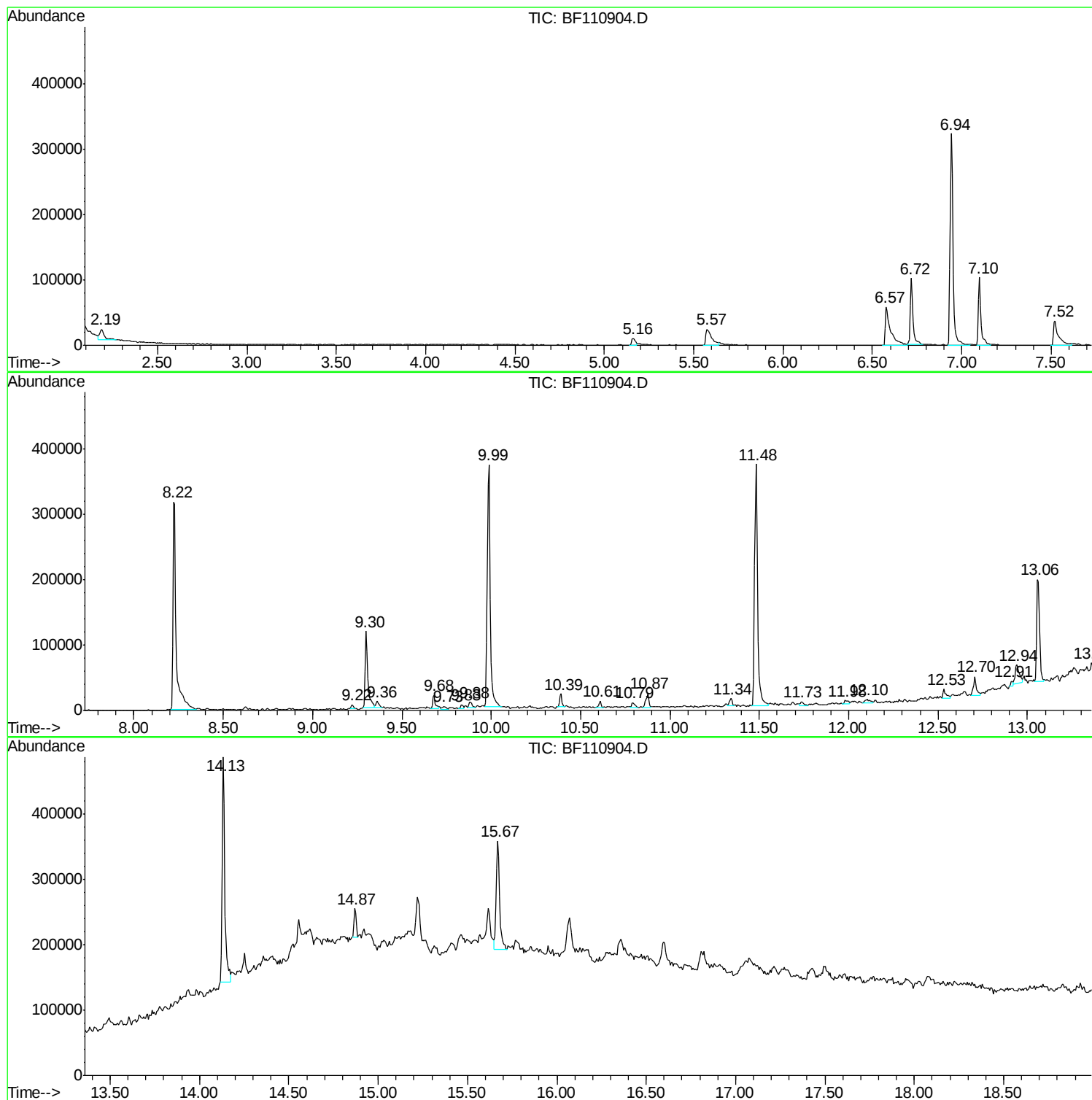
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BNA_F
ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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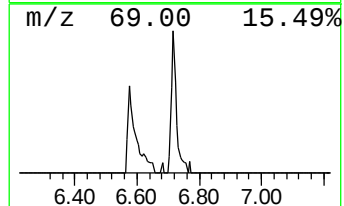
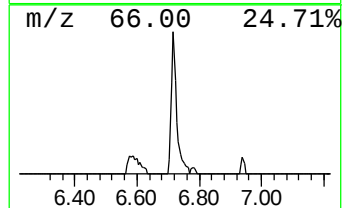
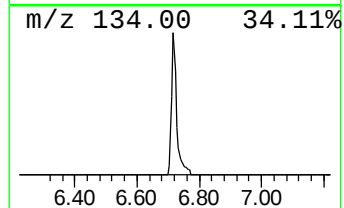
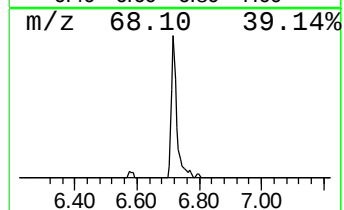
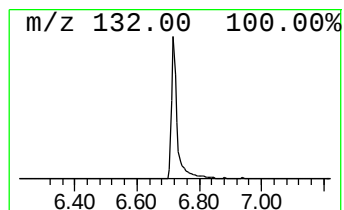
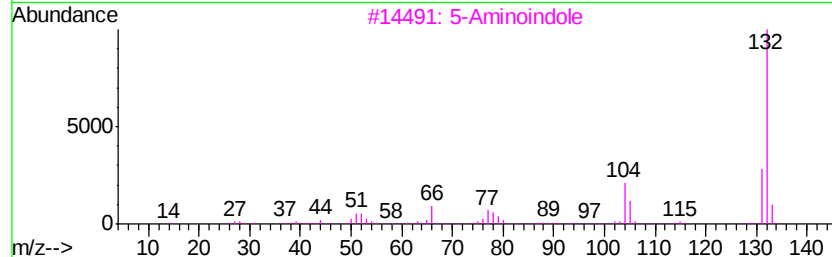
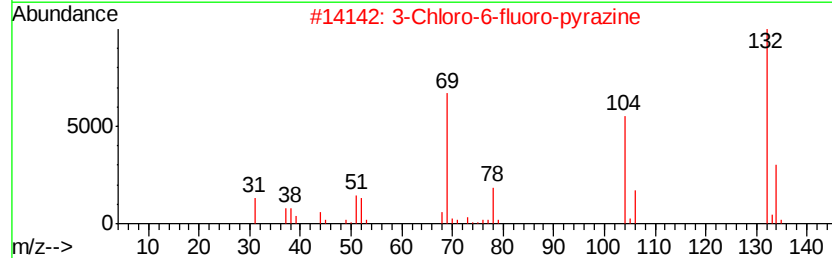
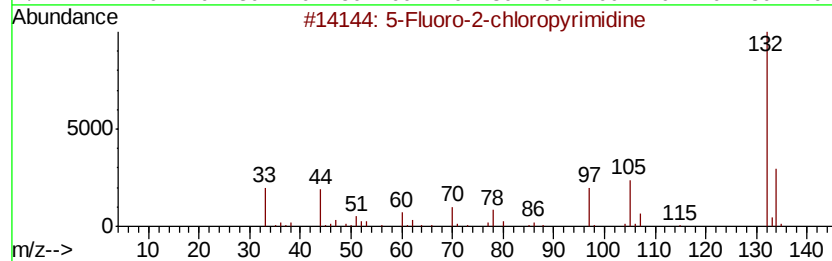
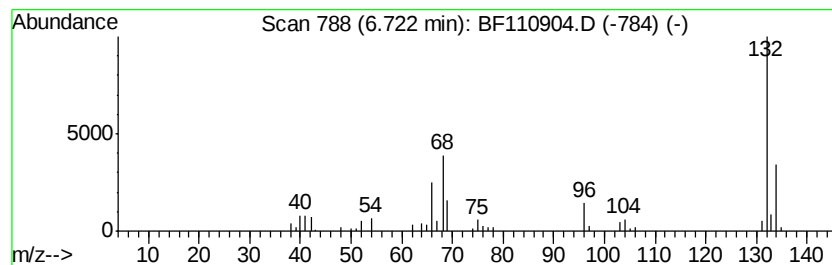
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.72 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	6.75 ng	112807	1,4-Dichlorobenzene-d4	6.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3			5-Aminoindole	132	C8H8N2	005192-03-0	10
4			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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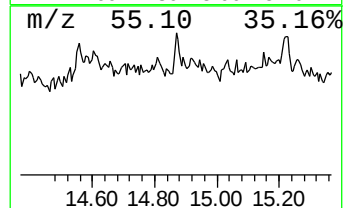
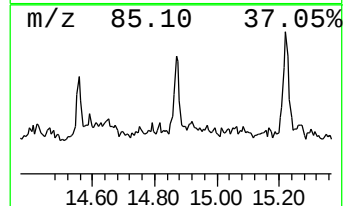
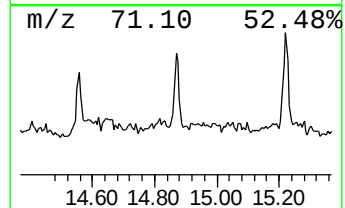
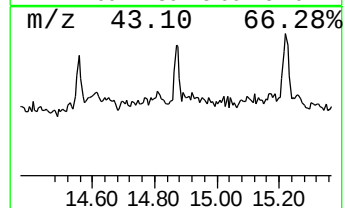
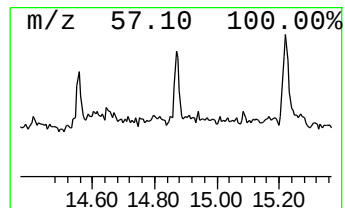
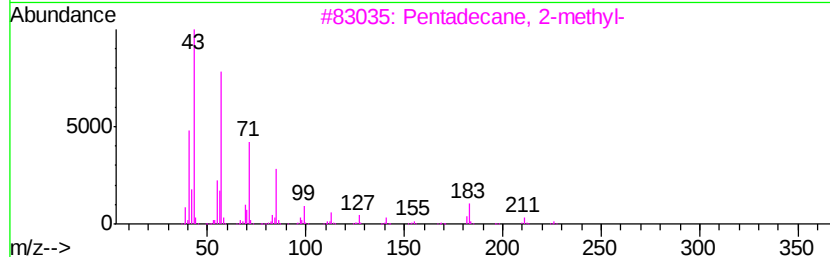
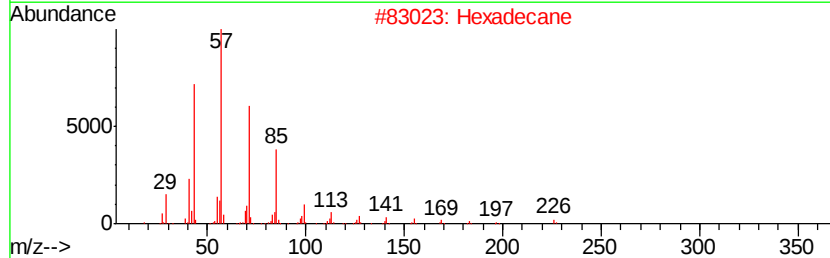
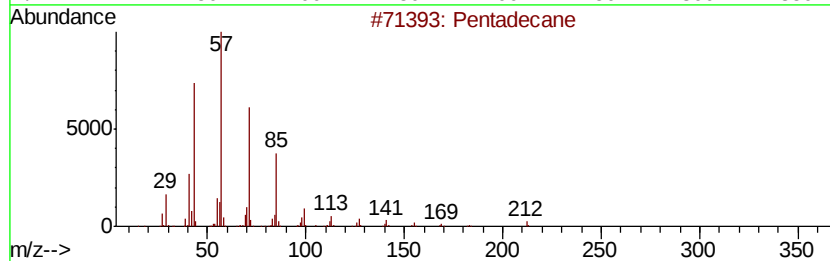
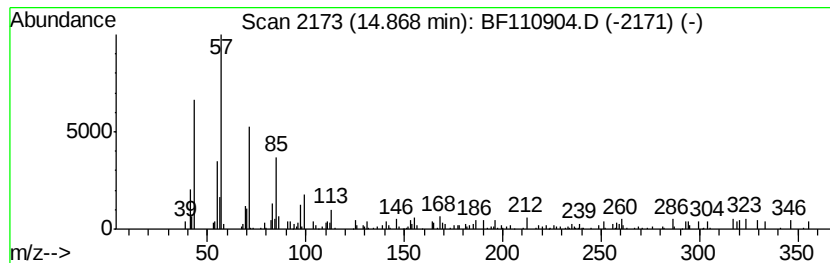
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Peak Number 4 Pentadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	2.22 ng	41144	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	60
2	Hexadecane		226	C16H34	000544-76-3	60
3	Pentadecane, 2-methyl-		226	C16H34	001560-93-6	53
4	Eicosane, 2-methyl-		296	C21H44	001560-84-5	53
5	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	49



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TIC Library : C:\DATABASE\NIST11.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.72	6.72	6.8	ng	112807	1	6.94	334477	20.0
Pentadecane	14.87	2.2	ng	41144	5	14.13	370913	20.0