

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071718\
 Data File : BF107450.D
 Acq On : 17 Jul 2018 15:37
 Operator : JU/SJ
 Sample : J3826-03 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-071618

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-BF071618.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	5.207	485	488	491	rBV3	32025	34751	2.85%	0.400%
2	5.601	550	555	559	rBV	278476	245686	20.15%	2.826%
3	6.601	722	725	733	rVB	285076	254742	20.89%	2.930%
4	6.754	748	751	756	rVB	318014	274640	22.52%	3.159%
5	6.984	787	790	794	rBV	909003	767975	62.97%	8.832%
6	7.142	813	817	820	rVB	244492	220159	18.05%	2.532%
7	7.542	882	885	890	rVB	190000	146924	12.05%	1.690%
8	8.272	1005	1009	1012	rBV	1075693	956268	78.41%	10.998%
9	9.342	1187	1191	1194	rBV	358646	302778	24.83%	3.482%
10	9.730	1254	1257	1261	rVB	42445	48193	3.95%	0.554%
11	10.030	1304	1308	1314	rVB2	1263964	1103876	90.52%	12.696%
12	10.819	1438	1442	1447	rBV2	245769	225539	18.49%	2.594%
13	10.983	1467	1470	1472	rBV4	17074	20451	1.68%	0.235%
14	11.366	1534	1535	1538	rBV3	24731	18201	1.49%	0.209%
15	11.401	1538	1541	1543	rBV4	21994	26202	2.15%	0.301%
16	11.460	1549	1551	1555	rBV5	14749	18437	1.51%	0.212%
17	11.524	1558	1562	1565	rVV	1353310	1219515	100.00%	14.026%
18	11.548	1565	1566	1570	rVB2	68824	43610	3.58%	0.502%
19	11.583	1570	1572	1574	rBV3	15406	18550	1.52%	0.213%
20	11.707	1591	1593	1595	rBV3	19718	19137	1.57%	0.220%
21	11.995	1640	1642	1645	rBV4	27900	31174	2.56%	0.359%
22	12.024	1645	1647	1650	rVV4	35145	42729	3.50%	0.491%
23	12.054	1650	1652	1654	rVB3	40075	32601	2.67%	0.375%
24	12.101	1657	1660	1663	rBV5	37844	63767	5.23%	0.733%
25	12.130	1663	1665	1669	rBV5	43907	61813	5.07%	0.711%
26	12.207	1675	1678	1679	rBV3	31119	27931	2.29%	0.321%
27	12.254	1684	1686	1692	rVV7	25923	50213	4.12%	0.577%
28	12.742	1766	1769	1771	rVB2	78830	74066	6.07%	0.852%
29	12.971	1805	1808	1812	rVB	105681	97956	8.03%	1.127%
30	13.107	1828	1831	1835	rBV	388630	378029	31.00%	4.348%
31	14.171	2009	2012	2016	rBV	1161426	1138337	93.34%	13.092%
32	15.718	2271	2275	2281	rVB	553839	730691	59.92%	8.404%

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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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071618.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

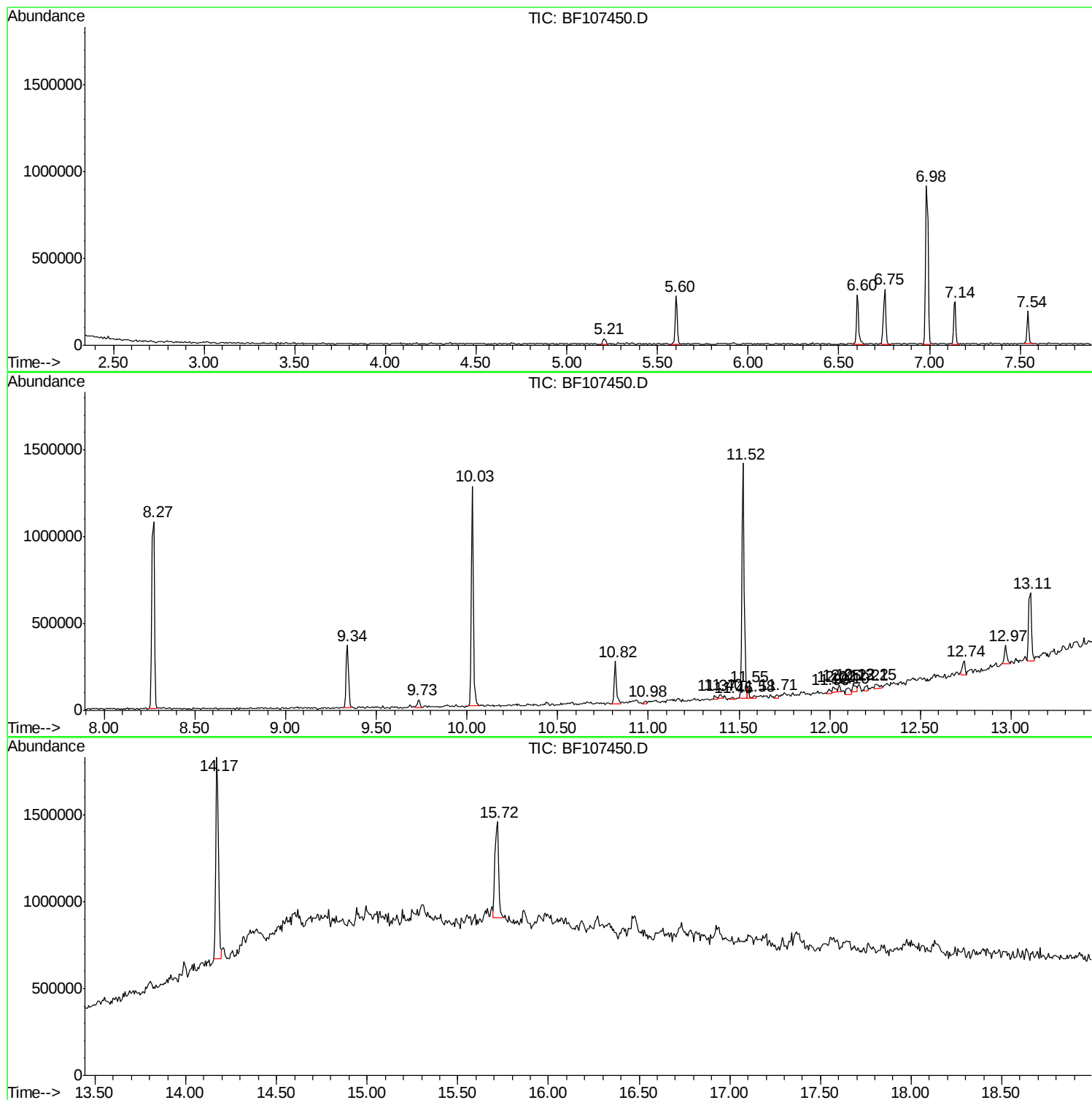
Sum of corrected areas: 8694941

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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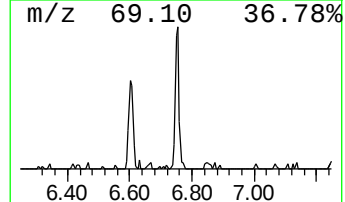
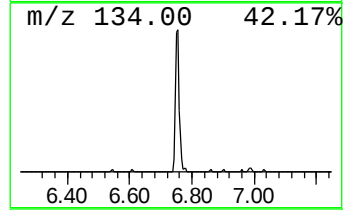
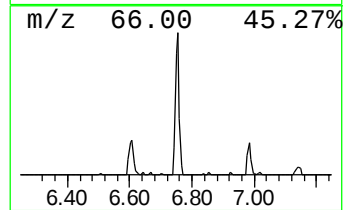
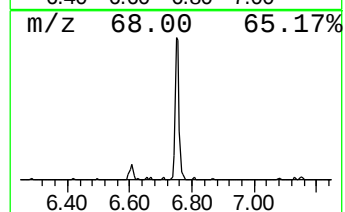
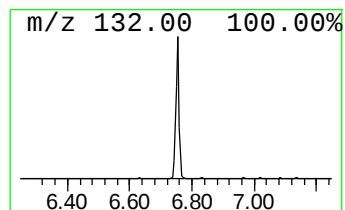
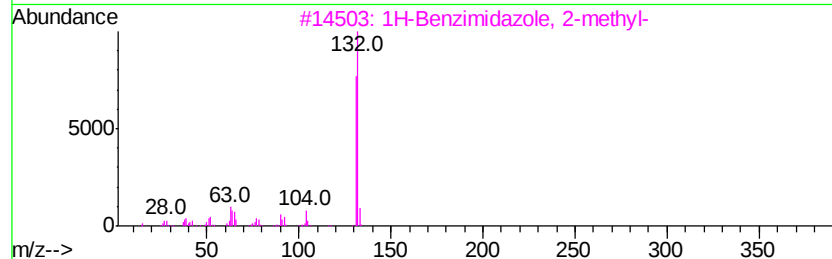
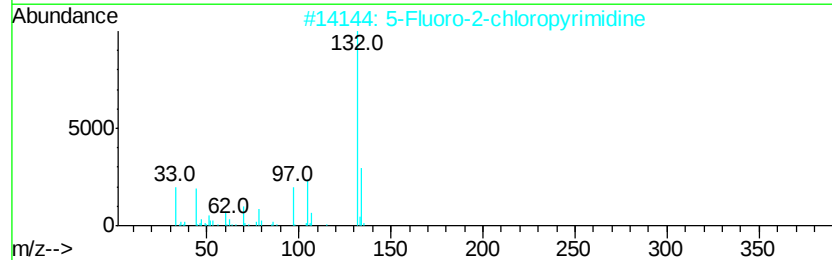
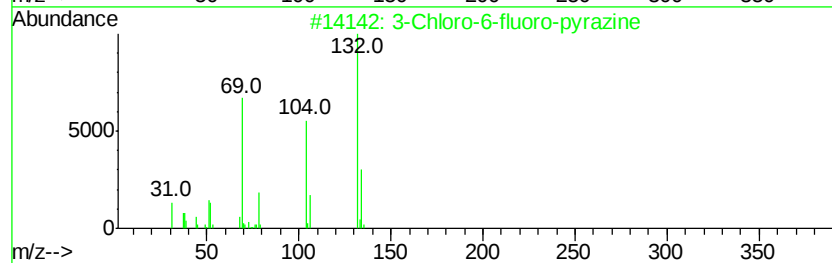
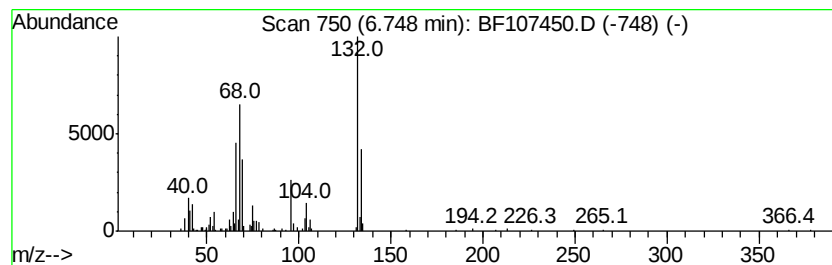
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	7.15 ng	274640	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	14
5			Benzeneethanamine, N-(2-phenylet...	223	C16H17N	005319-59-5	10



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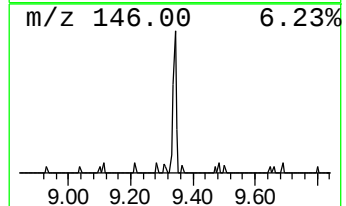
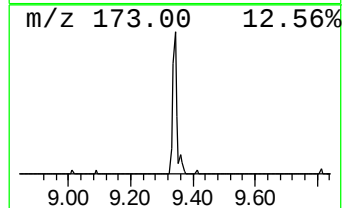
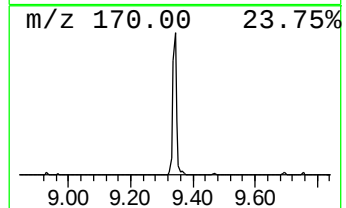
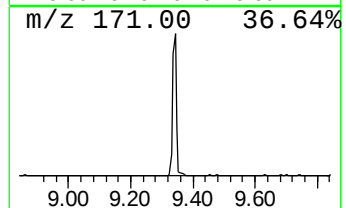
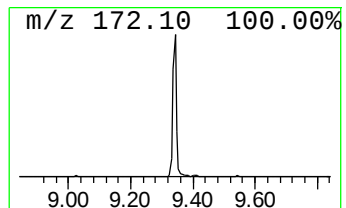
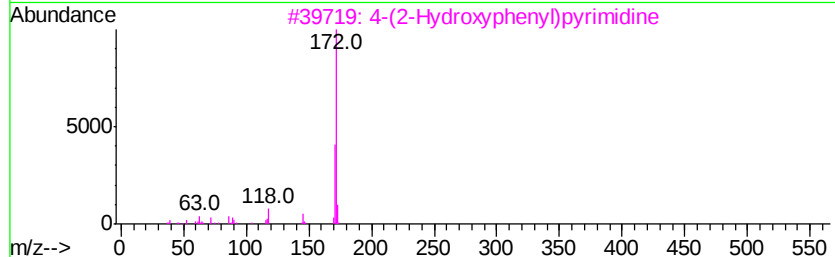
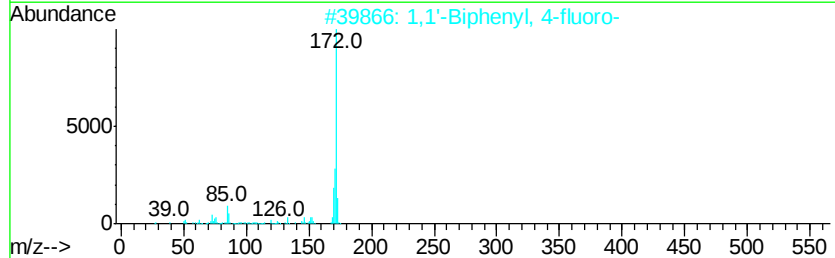
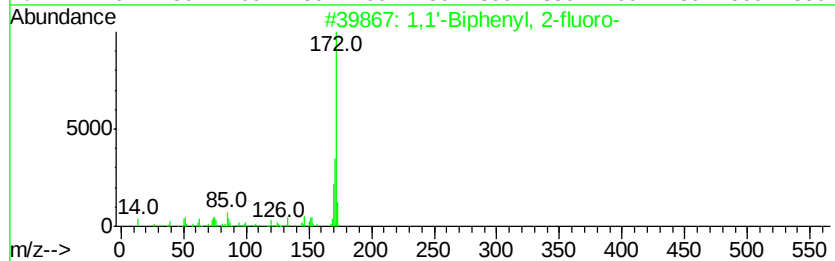
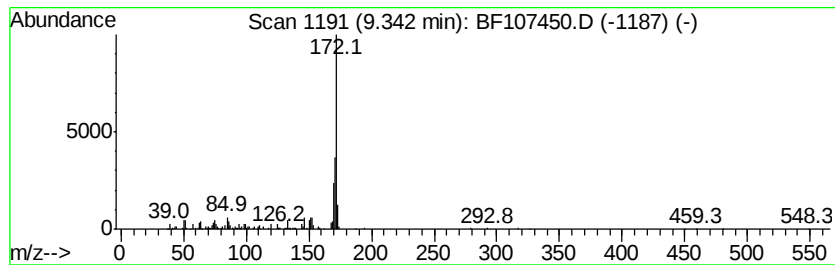
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Peak Number 3 1,1'-Biphenyl, 2-fluoro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.34	5.49 ng	302778	Acenaphthene-d10	10.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-fluoro-	172	C12H9F	000321-60-8	95
2	1,1'-Biphenyl, 4-fluoro-	172	C12H9F	000324-74-3	81
3	4-(2-Hydroxyphenyl)pyrimidine	172	C10H8N2O	068535-55-7	59
4	4-Pyrimidinol, 5-phenyl-	172	C10H8N2O	022433-69-8	53
5	D-Norleucine, N-propoxycarbonyl-...	273	C14H27NO4	1000338-42-8	47



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown6.75	6.75	7.2	ng	274640	1	6.98	767975	20.0
1,1'-Biphenyl, 2-...	9.34	5.5	ng	302778	3	10.03	1103880	20.0