

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052617\
 Data File : BF095481.D
 Acq On : 27 May 2017 3:21
 Operator : SJ/MA
 Sample : PB99316BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99316BL

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-BF050217.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	5.107	535	539	557	rBV2	61014	173358	7.12%	1.227%
2	5.725	640	644	680	rBV	436151	1320418	54.21%	9.345%
3	7.254	900	904	908	rBV	29795	45746	1.88%	0.324%
4	7.295	908	911	927	rVV	419120	1118175	45.91%	7.914%
5	7.413	927	931	957	rVB	873536	1686526	69.24%	11.937%
6	7.742	983	987	999	rBV	358941	460979	18.93%	3.263%
7	7.978	1023	1027	1047	rBV	1345341	1751585	71.91%	12.397%
8	8.648	1137	1141	1181	rBV	478324	1127059	46.27%	7.977%
9	9.772	1328	1332	1349	rBV	227083	547711	22.49%	3.877%
10	11.536	1627	1632	1647	rBV	1328085	2435657	100.00%	17.239%
11	12.607	1809	1814	1836	rBV2	215355	449652	18.46%	3.182%
12	13.918	2032	2037	2058	rBV	152597	394597	16.20%	2.793%
13	15.013	2218	2223	2232	rBV	185246	351331	14.42%	2.487%
14	17.318	2610	2615	2632	rBV	1240973	1669427	68.54%	11.816%
15	18.683	2843	2847	2869	rBV	180671	307361	12.62%	2.175%
16	20.365	3128	3133	3146	rBV2	148871	289416	11.88%	2.048%

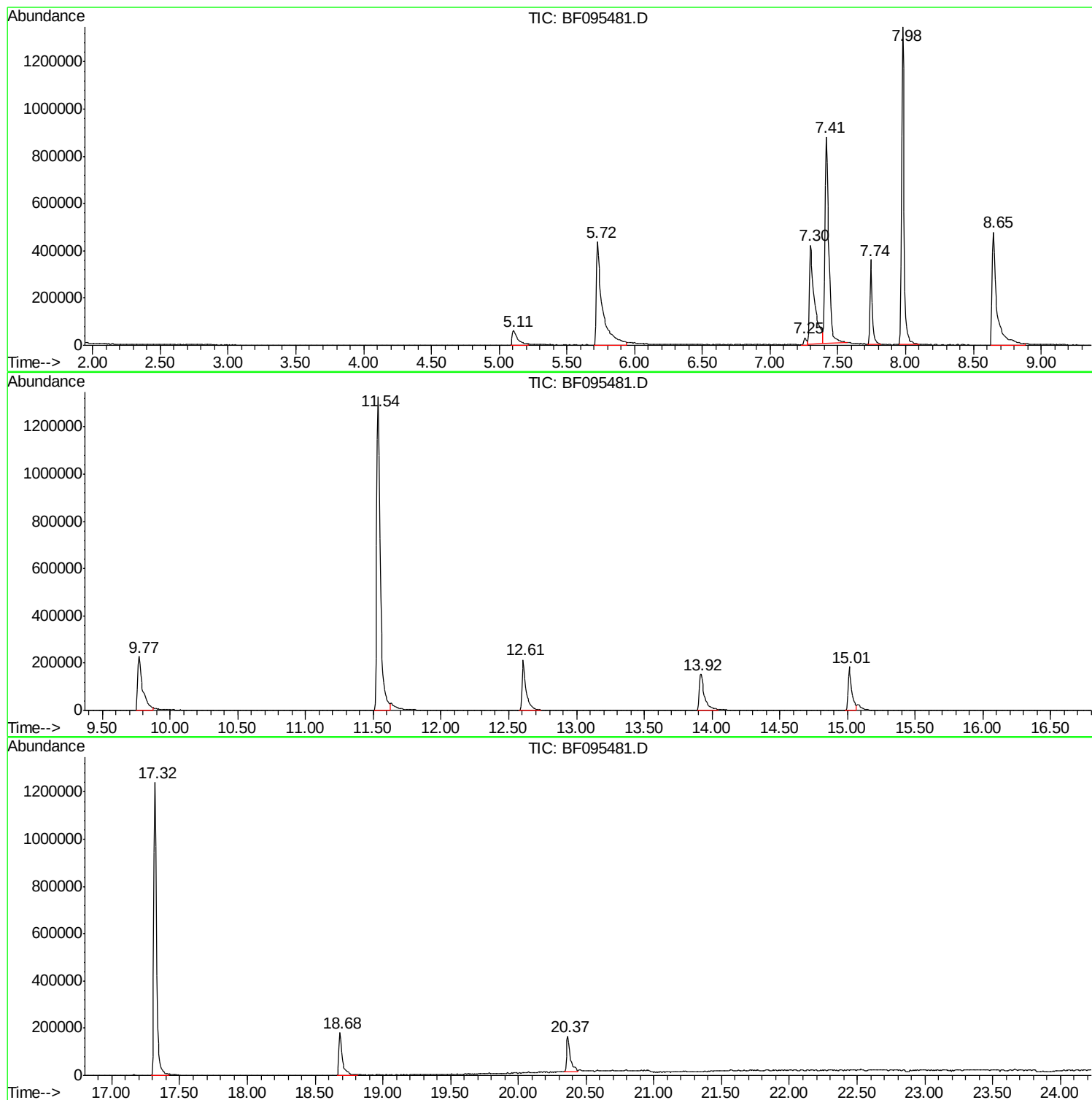
Sum of corrected areas: 14128998

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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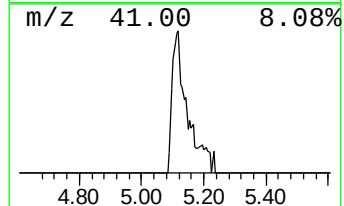
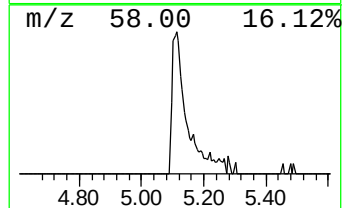
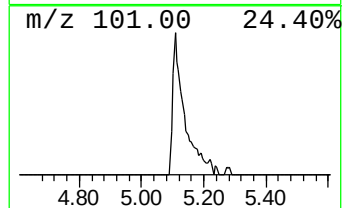
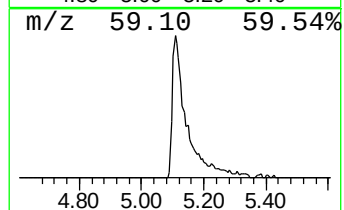
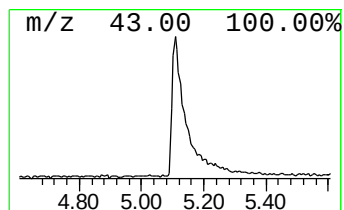
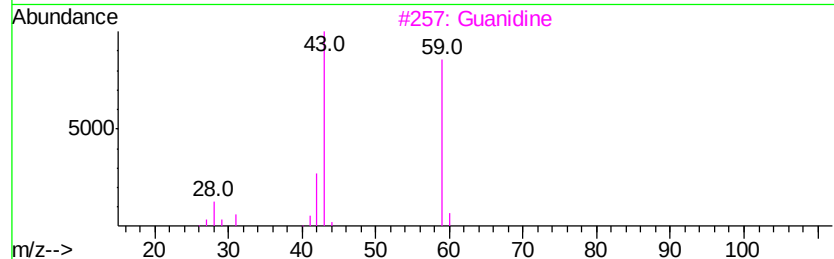
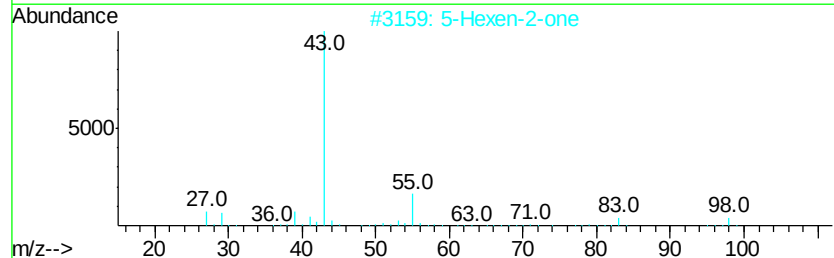
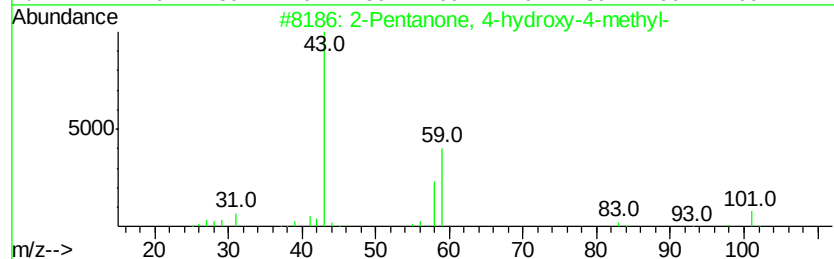
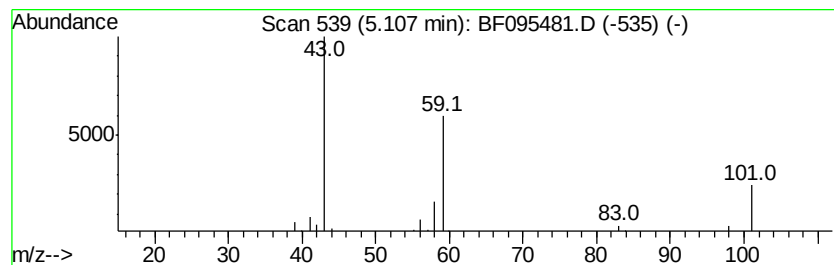
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	7.52 ng	173358	1,4-Dichlorobenzene-d4	7.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
3	Guanidine	59	CH5N3	000113-00-8	9		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9		
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9		



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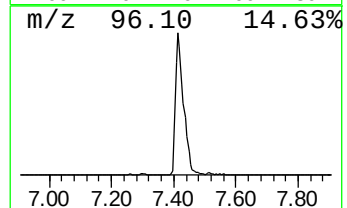
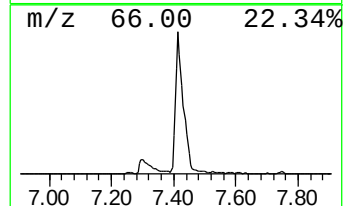
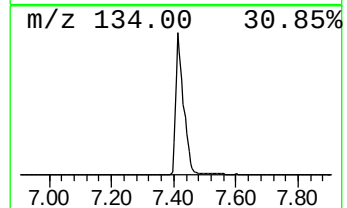
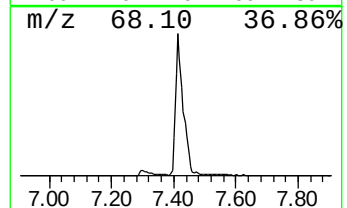
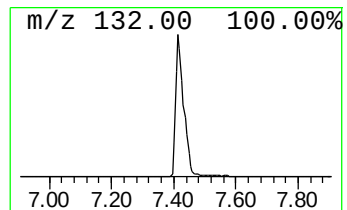
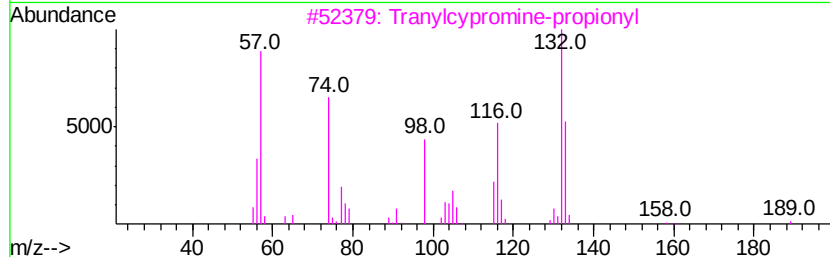
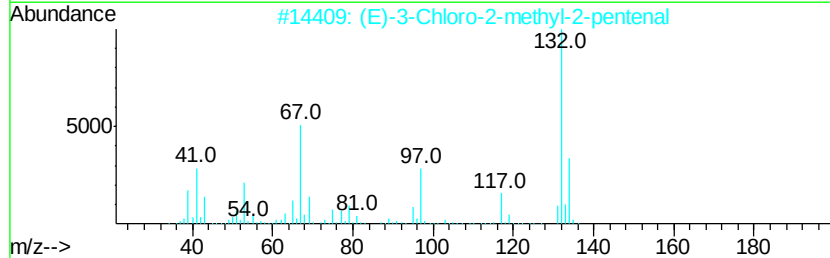
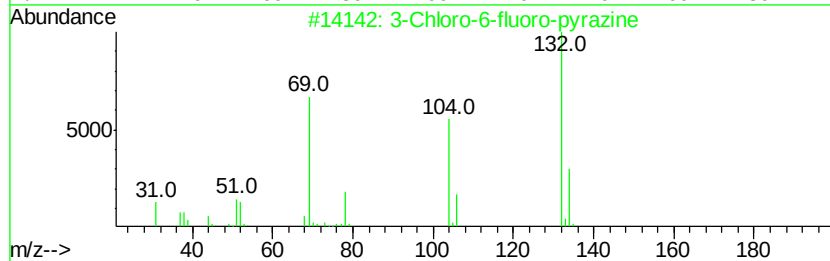
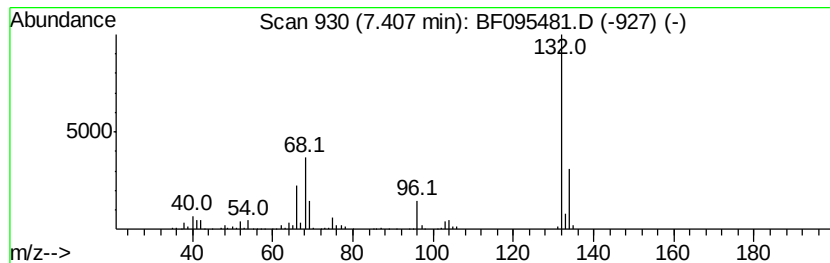
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Peak Number 2 unknown7.41 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.41	73.17 ng	1686530	1,4-Dichlorobenzene-d4	7.74

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	16
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.11	7.5	ng	173358	1	7.74	460979	20.0
unknown7.41	7.41	73.2	ng	1686530	1	7.74	460979	20.0