

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092518\
 Data File : BF109298.D
 Acq On : 25 Sep 2018 20:49
 Operator : JU/SJ
 Sample : J5070-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-RBR-200028

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-BF092218.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	4.910	476	480	490	rBV	176310	243519	7.83%	1.035%
2	5.328	546	551	555	rBV	2319143	2444255	78.60%	10.390%
3	6.351	720	725	728	rBV	2434625	2307390	74.19%	9.808%
4	6.481	743	747	750	rBV	2843017	2770191	89.08%	11.775%
5	6.710	782	786	789	rBV	904596	687589	22.11%	2.923%
6	6.869	809	813	816	rBV	2693310	2250625	72.37%	9.567%
7	7.269	876	881	884	rBV	1762391	1786189	57.44%	7.592%
8	7.992	1000	1004	1006	rBV	992861	851755	27.39%	3.620%
9	8.010	1006	1007	1011	rVB	112737	68306	2.20%	0.290%
10	9.069	1182	1187	1190	rBV	3672073	3109916	100.00%	13.219%
11	9.739	1297	1301	1305	rBV2	1038268	847754	27.26%	3.603%
12	10.522	1430	1434	1452	rBV	1513150	1529639	49.19%	6.502%
13	11.216	1548	1552	1556	rBV	911952	747890	24.05%	3.179%
14	12.798	1817	1821	1825	rBV	2671582	2628181	84.51%	11.171%
15	13.845	1995	1999	2002	rBV	810273	732966	23.57%	3.116%
16	15.245	2233	2237	2242	rBV	467624	519792	16.71%	2.209%

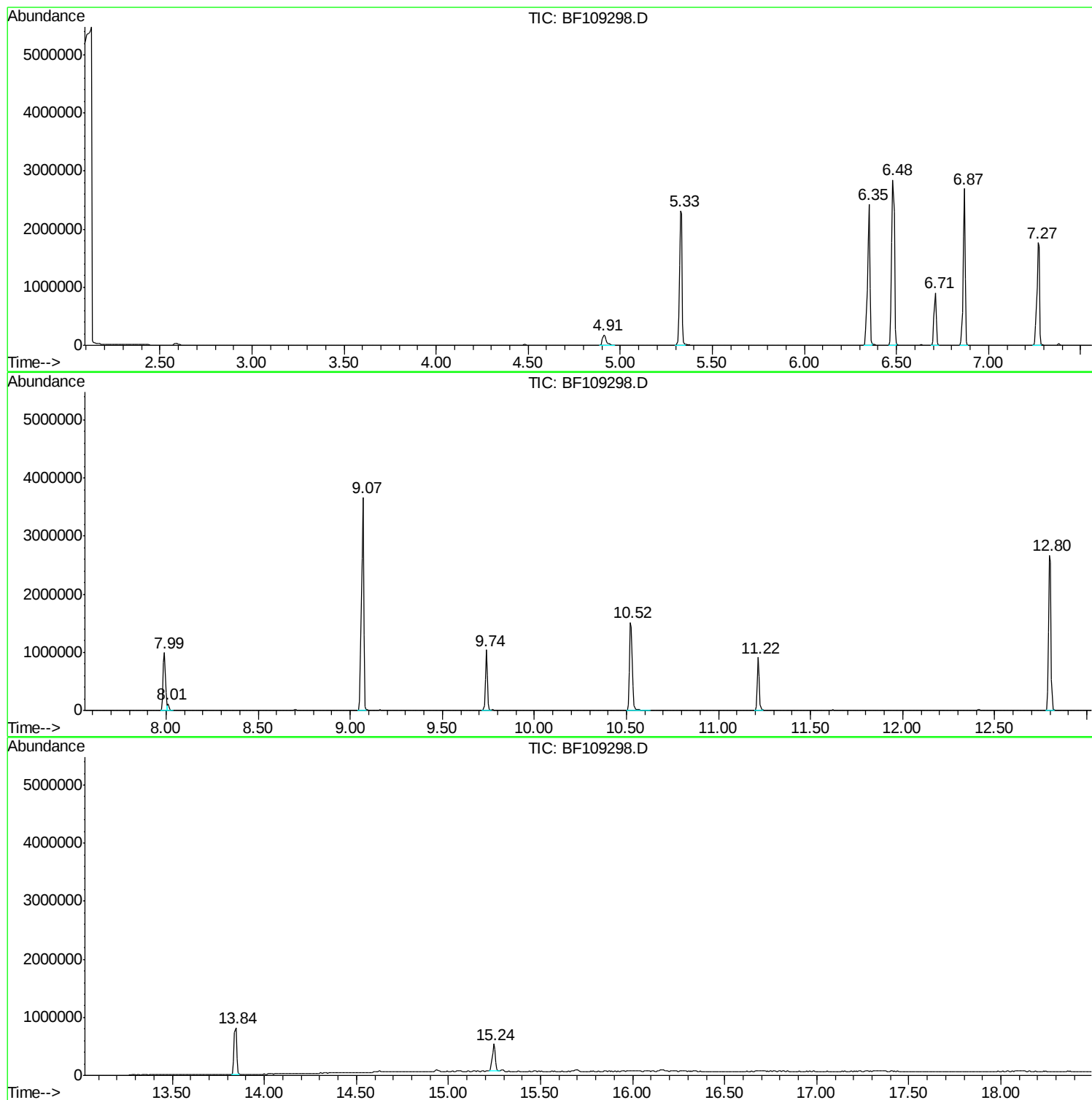
Sum of corrected areas: 23525957

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.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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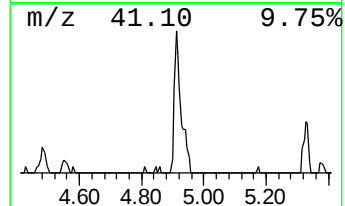
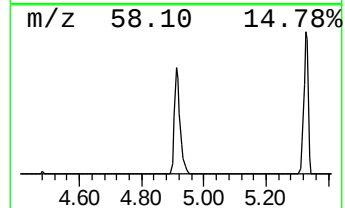
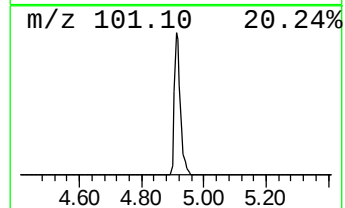
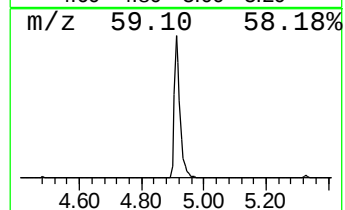
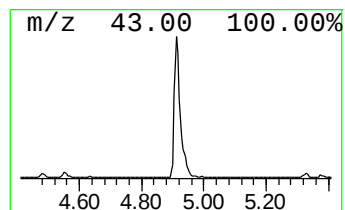
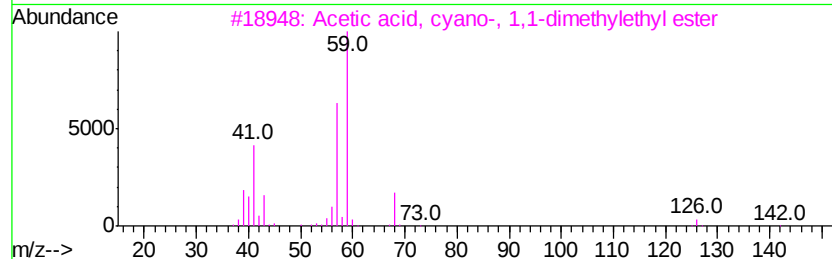
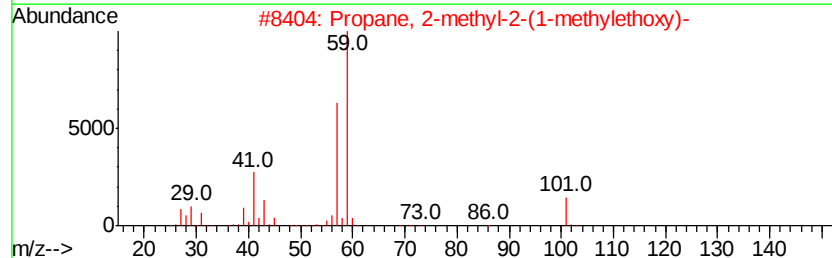
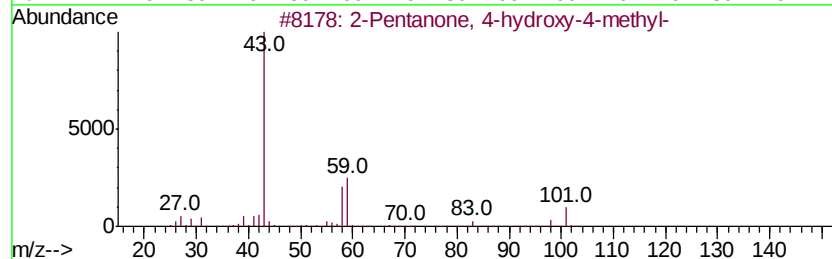
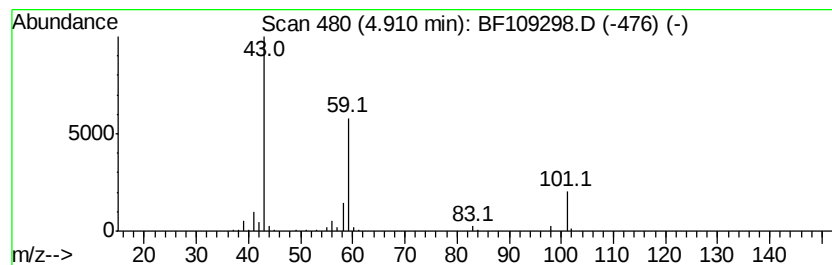
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	7.08 ng	243519	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	28
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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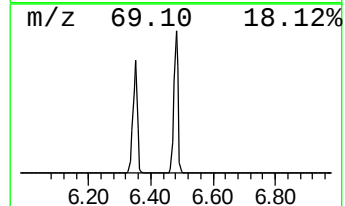
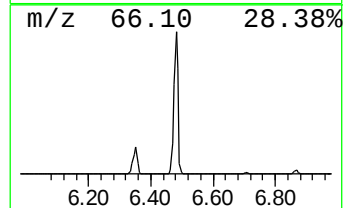
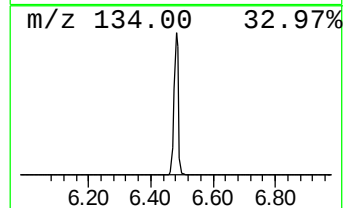
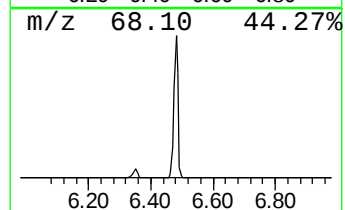
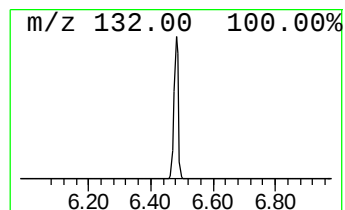
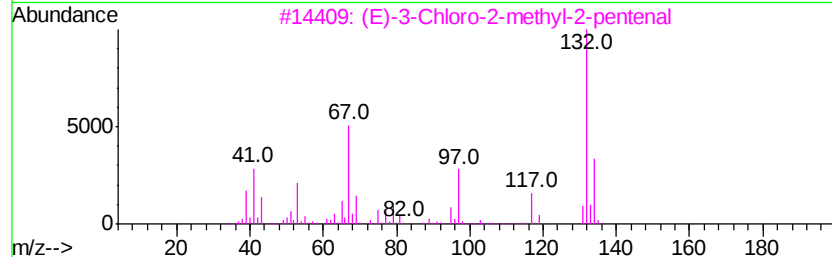
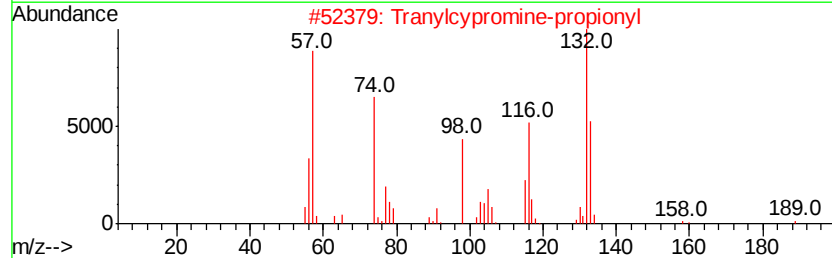
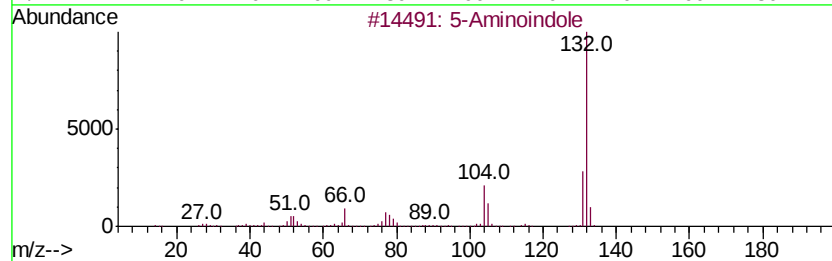
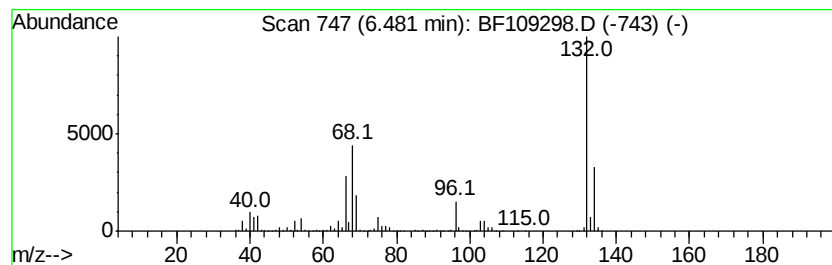
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Peak Number 2 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	80.58 ng	2770190	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Aminoindole	132	C8H8N2	005192-03-0	12
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	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...	4.91	7.1	ng	243519	1	6.71	687589	20.0
unknown6.48	6.48	80.6	ng	2770190	1	6.71	687589	20.0