

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
 Data File : BM015792.D
 Acq On : 06 Jul 2018 08:57
 Operator : SJ/JU
 Sample : PB110755BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110755BL

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_M\METHODS\8270-BM070518.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	5.328	342	347	363	rBV	278656	438572	6.02%	0.962%
2	5.810	423	429	447	rBV	2905781	4398524	60.42%	9.645%
3	7.463	703	710	724	rBV	2913542	4656589	63.96%	10.211%
4	7.834	766	773	786	rBV	3112379	4951537	68.01%	10.858%
5	8.310	847	854	870	rBV	517342	838106	11.51%	1.838%
6	8.634	900	909	926	rBV	2221819	3620039	49.73%	7.938%
7	9.498	1049	1056	1071	rBV	1687372	2831473	38.89%	6.209%
8	11.139	1328	1335	1343	rBV	605606	1026642	14.10%	2.251%
9	13.557	1739	1746	1756	rBV	4289136	6118731	84.05%	13.417%
10	14.927	1973	1979	1989	rBV2	846767	1267040	17.40%	2.778%
11	16.410	2224	2231	2247	rBV	3491385	4762763	65.42%	10.444%
12	17.656	2437	2443	2452	rBV	969102	1416862	19.46%	3.107%
13	20.215	2872	2878	2884	rBV	6361910	7280075	100.00%	15.964%
14	21.762	3136	3141	3149	rBV	863056	1136850	15.62%	2.493%
15	24.168	3544	3550	3560	rVB	423864	859433	11.81%	1.885%

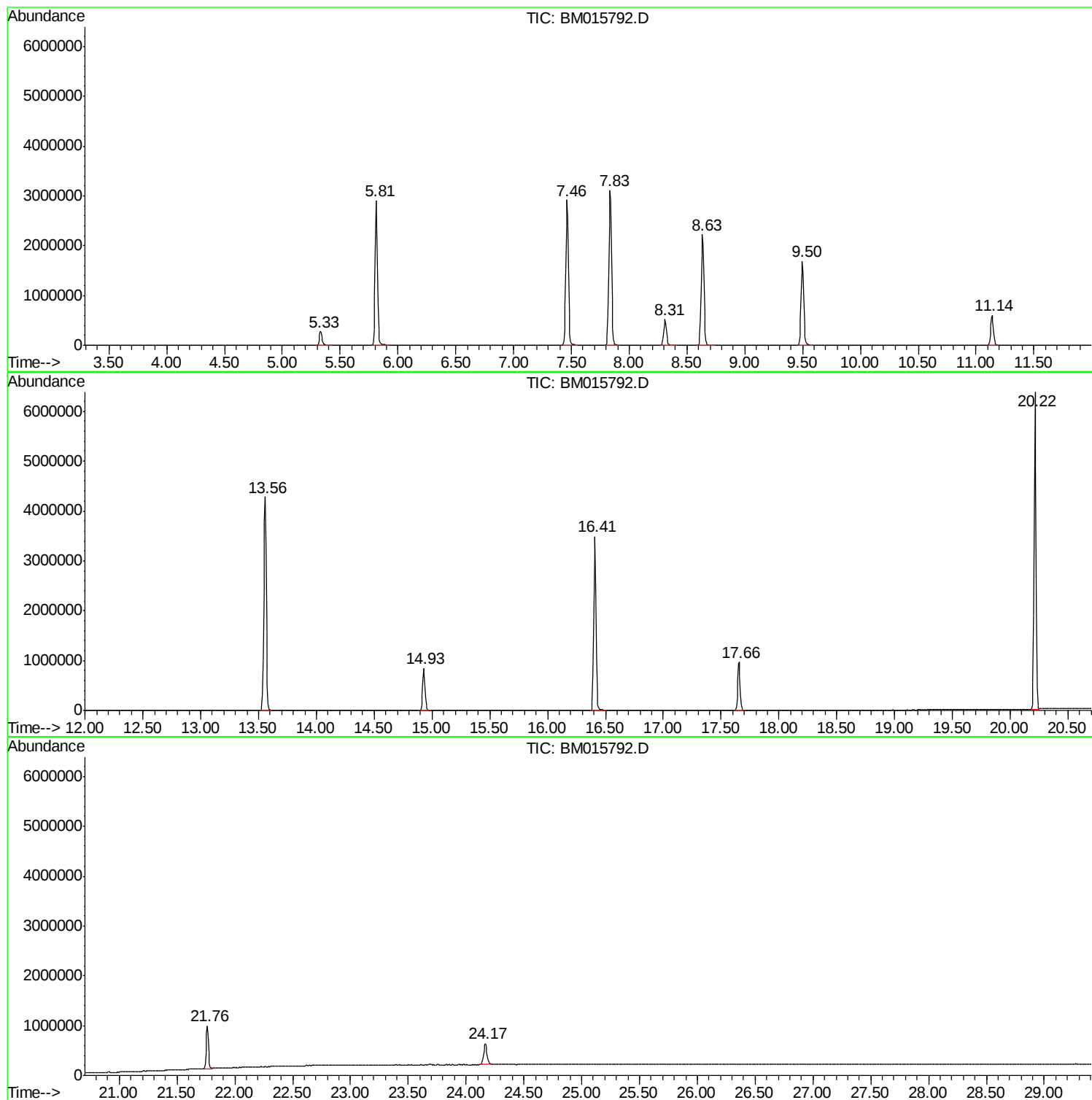
Sum of corrected areas: 45603236

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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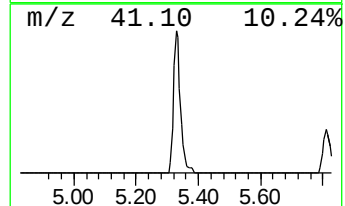
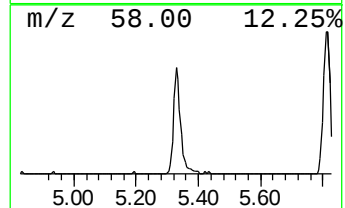
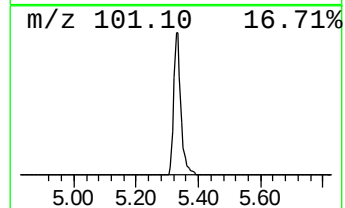
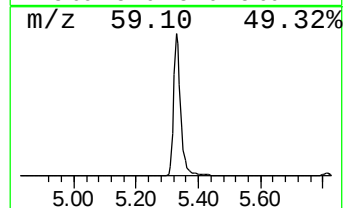
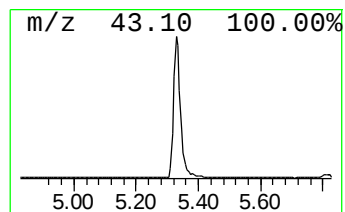
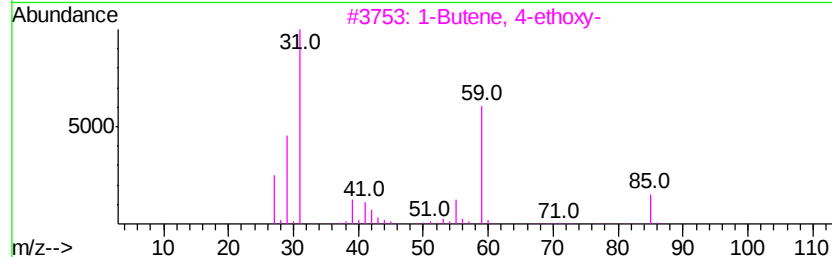
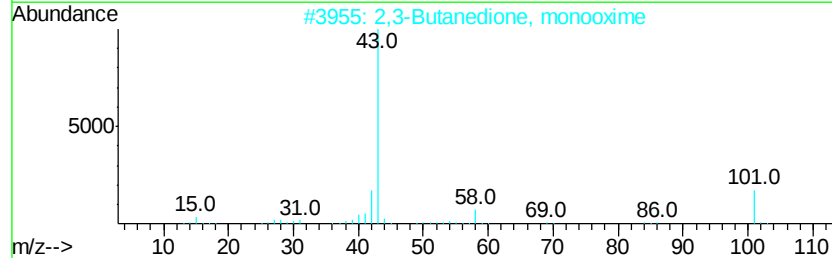
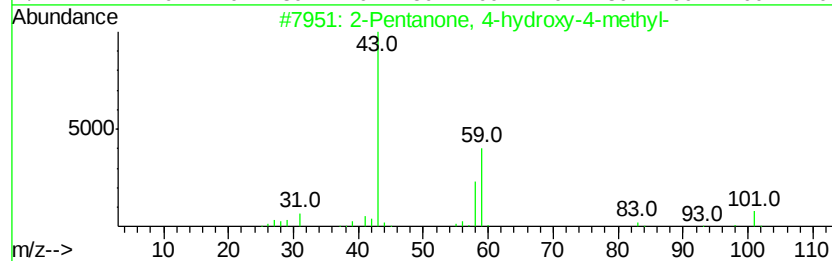
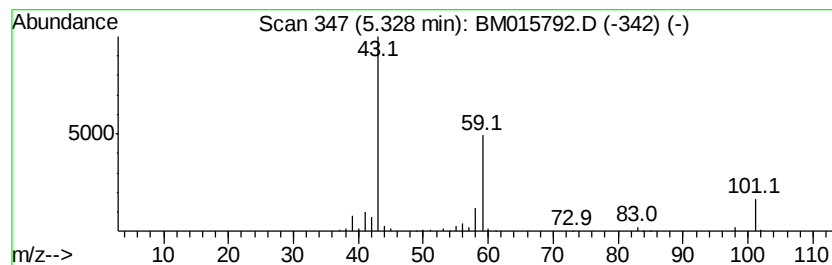
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.33	10.47 ng	438572	1,4-Dichlorobenzene-d4	8.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
3			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
4			Butane	58	C4H10	000106-97-8	9
5			Propylamine	59	C3H9N	000107-10-8	9



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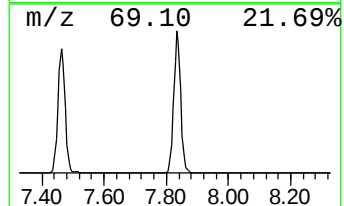
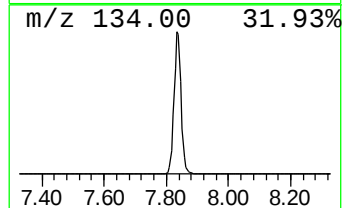
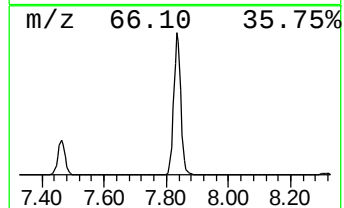
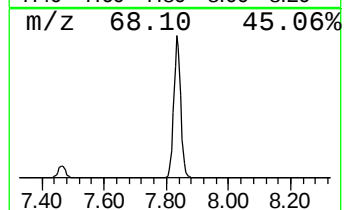
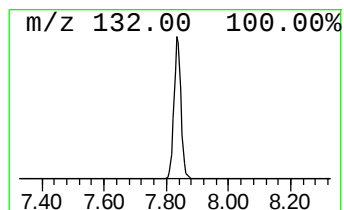
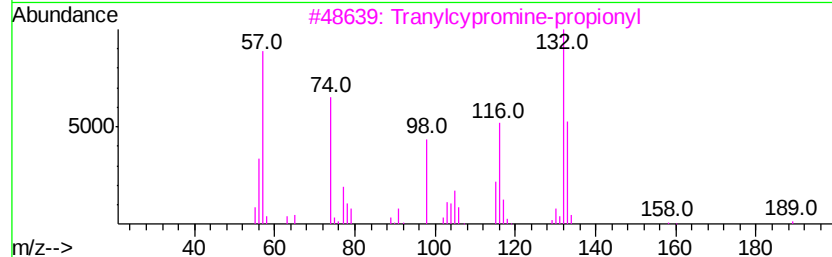
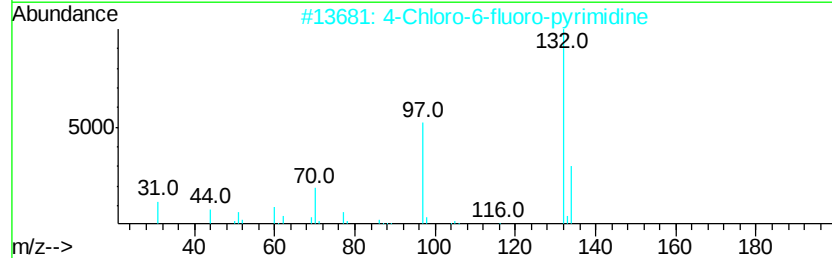
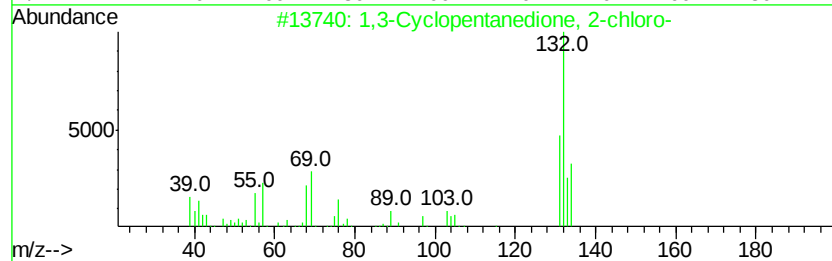
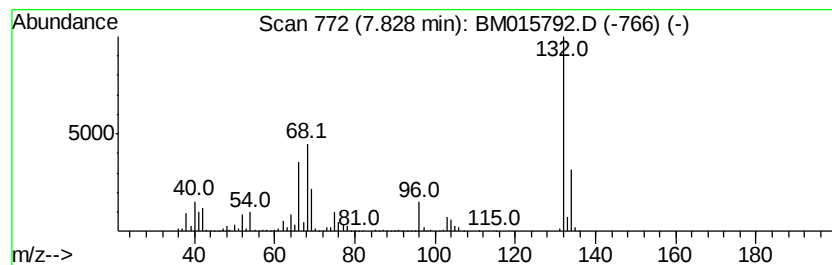
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Peak Number 2 unknown7.83 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	118.16 ng	4951540	1,4-Dichlorobenzene-d4	8.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.33	10.5	ng	438572	1	8.31	838106	20.0
unknown7.83	7.83	118.2	ng	4951540	1	8.31	838106	20.0