

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036642.D
 Acq On : 11 Sep 2018 2:09
 Operator : JU/SJ
 Sample : J4830-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090618-DBRI-E-D-M

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_G\METHODS\8270-BG082318.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	5.155	313	318	328	rVB	38785	59725	1.27%	0.310%
2	5.625	391	398	407	rBV	340855	563019	11.93%	2.923%
3	7.206	661	667	679	rBV2	236812	411658	8.72%	2.137%
4	7.582	724	731	740	rBV	635154	1088748	23.08%	5.653%
5	8.052	804	811	818	rBV	195537	334213	7.08%	1.735%
6	8.375	859	866	876	rVB	832286	1450636	30.75%	7.532%
7	9.215	1000	1009	1018	rBV	740856	1320729	27.99%	6.857%
8	10.860	1279	1289	1298	rVB	274506	501385	10.63%	2.603%
9	13.305	1697	1705	1712	rBV	2085419	3241713	68.71%	16.831%
10	14.121	1839	1844	1852	rBV	168802	252387	5.35%	1.310%
11	14.679	1932	1939	1947	rBV2	464450	758275	16.07%	3.937%
12	15.966	2153	2158	2163	rBV	43593	57152	1.21%	0.297%
13	16.160	2184	2191	2212	rBV	945321	1523401	32.29%	7.910%
14	17.423	2399	2406	2415	rBV	602149	958739	20.32%	4.978%
15	20.026	2843	2849	2861	rBV	3343885	4718216	100.00%	24.497%
16	21.683	3124	3131	3150	rBV2	604532	1066071	22.59%	5.535%
17	23.375	3413	3419	3424	rBV	30510	57118	1.21%	0.297%
18	24.897	3668	3678	3695	rBV2	264080	897162	19.01%	4.658%

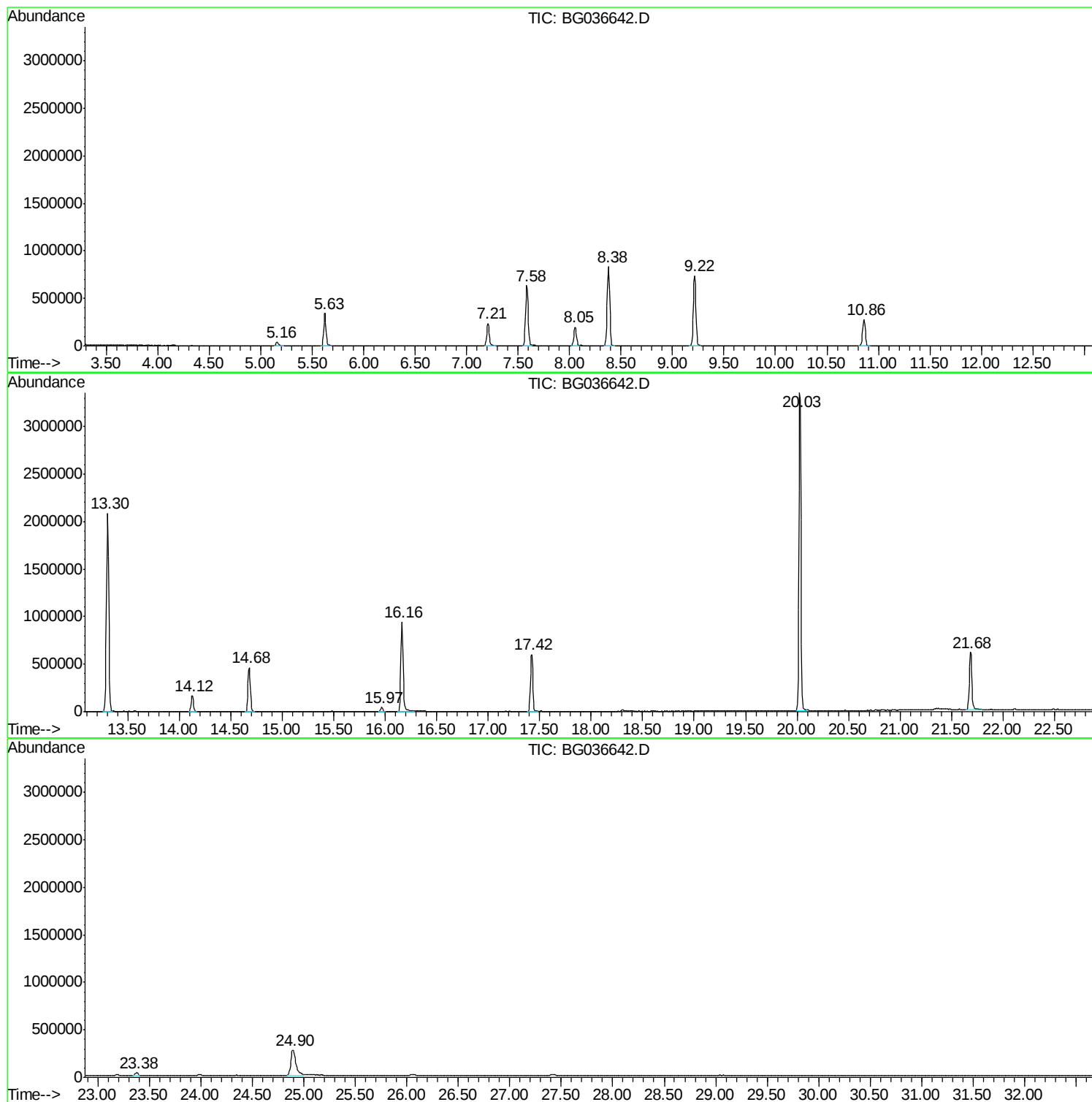
Sum of corrected areas: 19260347

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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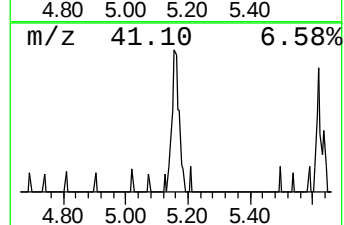
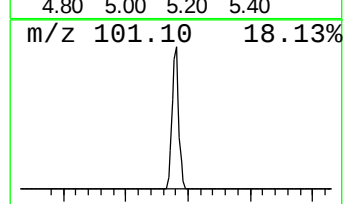
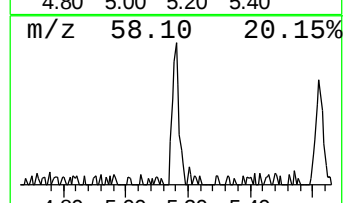
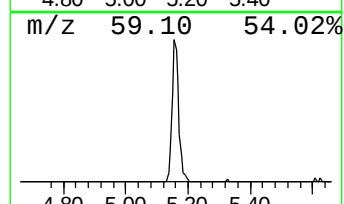
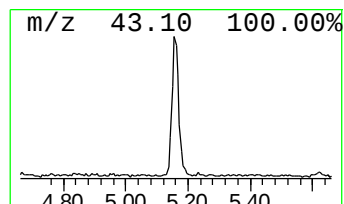
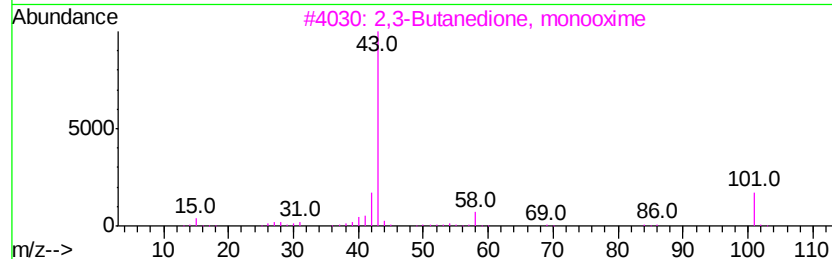
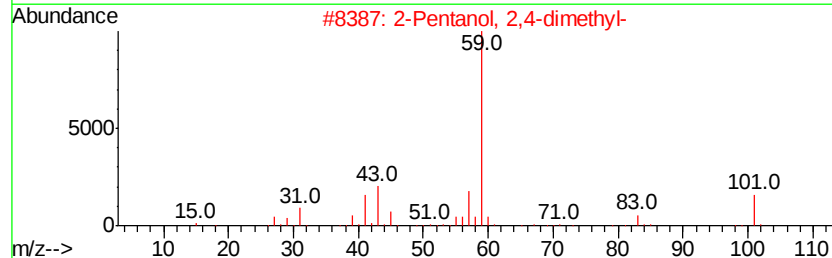
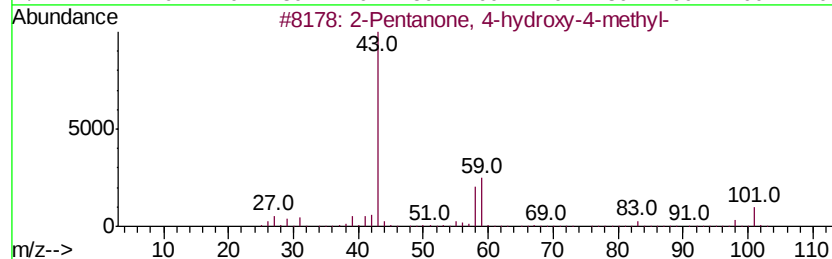
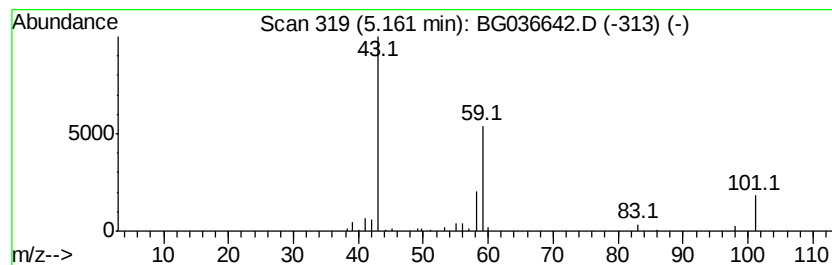
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	3.57 ng	59725	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			3-Heptanol	116	C7H16O	000589-82-2	25
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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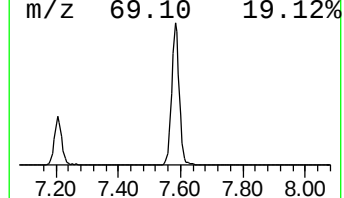
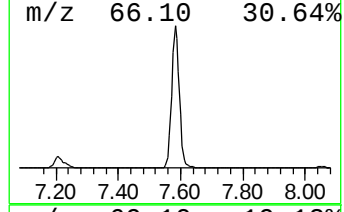
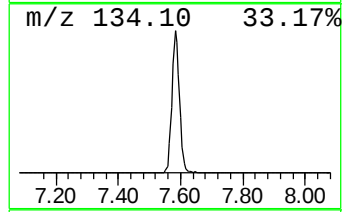
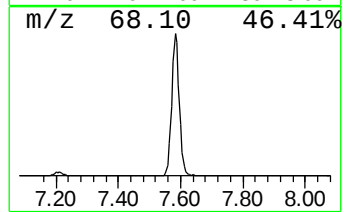
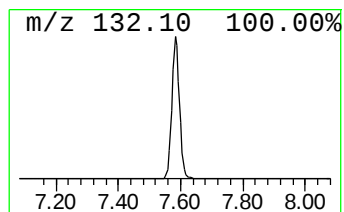
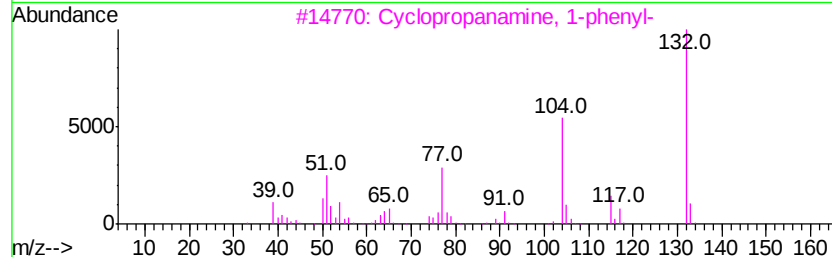
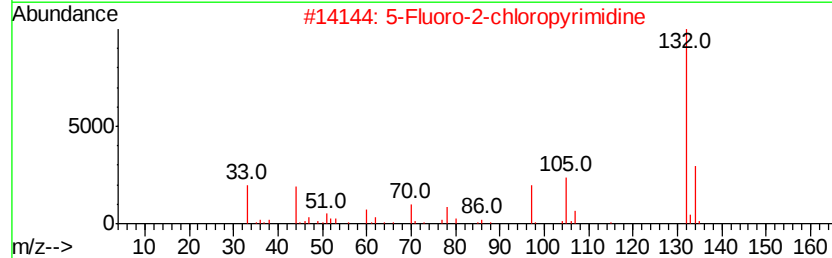
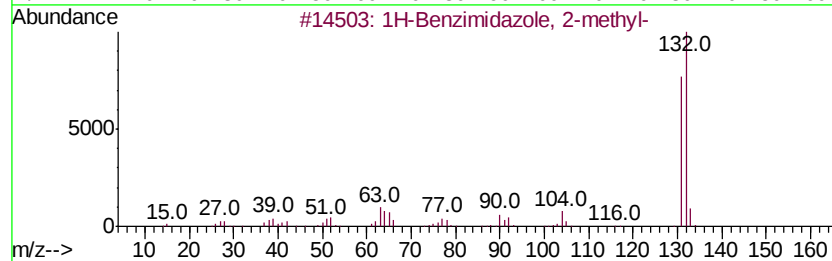
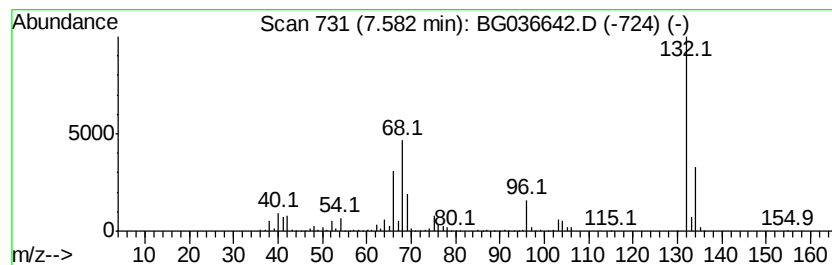
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	65.15 ng	1088750	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	3.6	ng	59725	1	8.05	334213	20.0
unknown7.58	7.58	65.2	ng	1088750	1	8.05	334213	20.0