

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110518\
 Data File : BG037808.D
 Acq On : 5 Nov 2018 18:26
 Operator : JU/SJ
 Sample : J5769-22
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FB-10-31-18

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_G\METHODS\8270-BG101818.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	5.174	315	321	329	rBV2	32996	56633	1.43%	0.310%
2	5.639	392	400	413	rBV	437586	745687	18.80%	4.077%
3	7.237	665	672	683	rBV	301227	526382	13.27%	2.878%
4	7.619	729	737	747	rBV	784885	1407927	35.49%	7.698%
5	8.095	811	818	830	rBV	193089	340692	8.59%	1.863%
6	8.418	865	873	883	rVB	737237	1308599	32.99%	7.155%
7	9.276	1010	1019	1033	rBV	609527	1175523	29.63%	6.427%
8	10.915	1290	1298	1310	rBV	283900	536465	13.52%	2.933%
9	13.353	1705	1713	1724	rBV	1815783	2977777	75.07%	16.281%
10	14.728	1940	1947	1959	rBV2	488866	799691	20.16%	4.372%
11	16.214	2194	2200	2220	rBV	1140924	1867425	47.08%	10.210%
12	17.478	2406	2415	2427	rBV	572577	921635	23.23%	5.039%
13	20.075	2850	2857	2871	rBV	2848610	3966853	100.00%	21.689%
14	21.761	3136	3144	3158	rBV	472708	950830	23.97%	5.199%
15	25.098	3696	3712	3727	rBV3	170023	707621	17.84%	3.869%

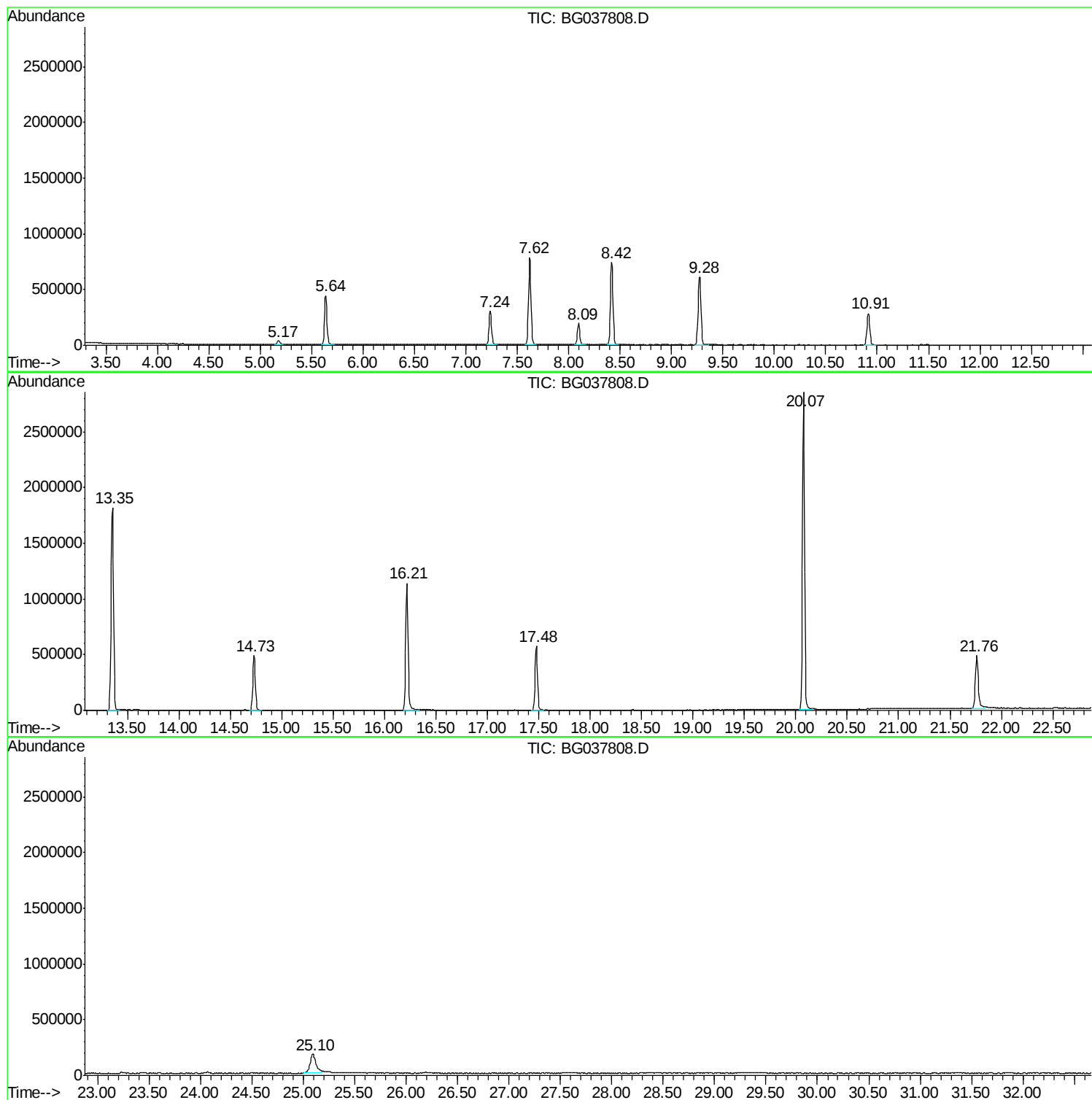
Sum of corrected areas: 18289740

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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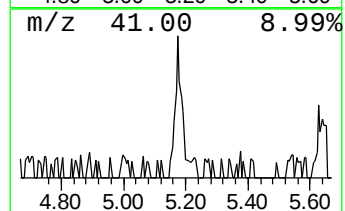
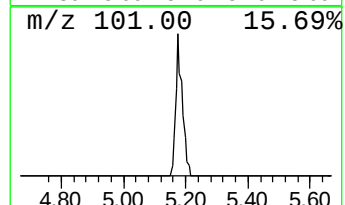
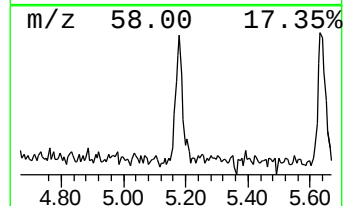
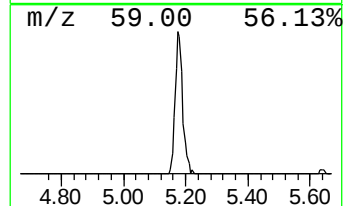
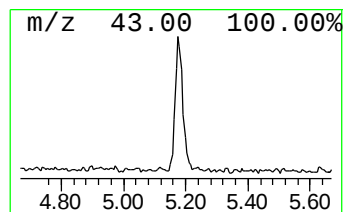
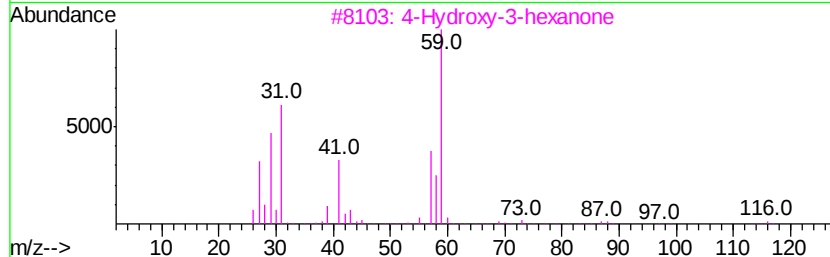
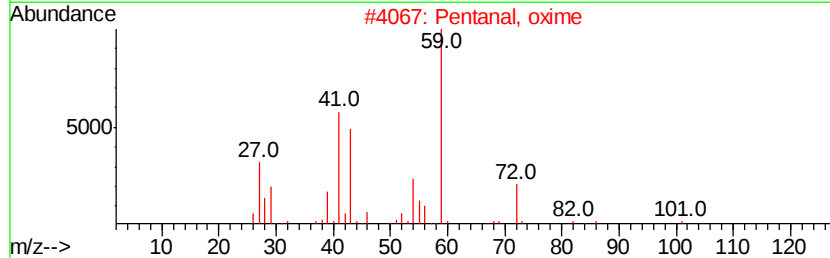
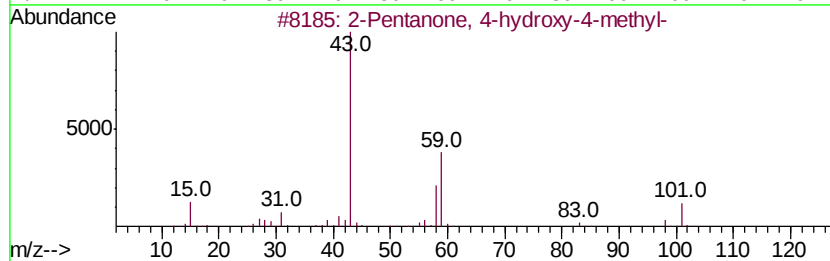
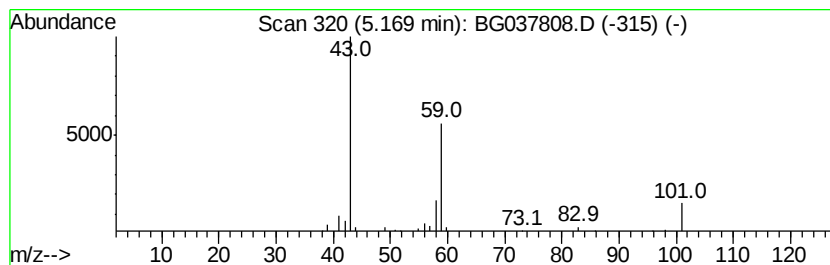
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.32 ng	56633	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Pentanal, oxime	101	C5H11NO	000628-79-5	25
3		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
4		2,3-Butanediol, 2,3-dimethyl-	118	C6H14O2	000076-09-5	17
5		1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	9



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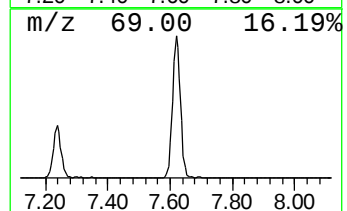
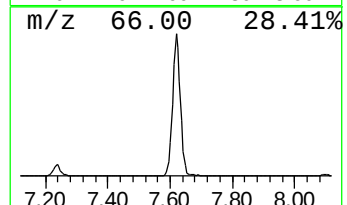
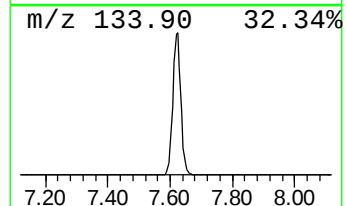
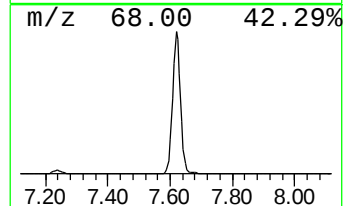
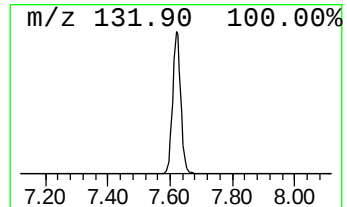
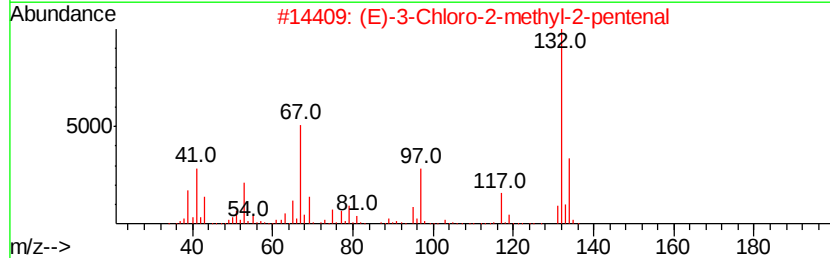
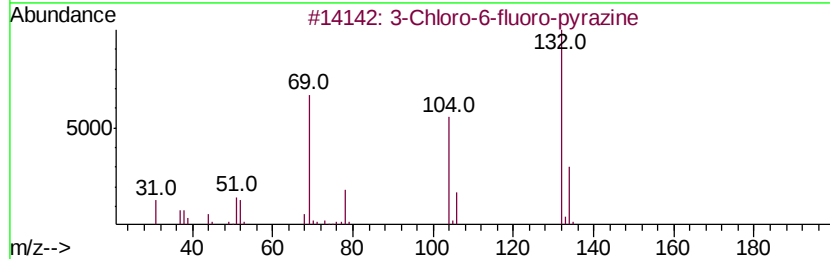
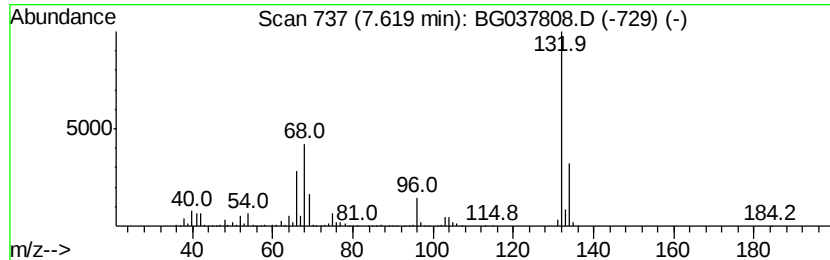
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	82.65 ng	1407930	1,4-Dichlorobenzene-d4	8.09

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	3.3	ng	56633	1	8.09	340692	20.0
unknown7.62	7.62	82.7	ng	1407930	1	8.09	340692	20.0