

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
 Data File : BG031566.D
 Acq On : 14 Dec 2017 19:48
 Operator : SJ/JU
 Sample : I6883-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FB-02

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_G\METHODS\8270-BG121117.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	5.175	315	321	331	rBV	63124	104338	1.48%	0.401%
2	5.657	397	403	418	rBV	453710	830709	11.80%	3.192%
3	7.266	671	677	692	rBV	268035	528223	7.50%	2.030%
4	7.631	732	739	758	rBV	935104	1691172	24.01%	6.499%
5	8.095	809	818	831	rBV	231190	414907	5.89%	1.594%
6	8.424	866	874	886	rBV	913130	1664741	23.64%	6.397%
7	9.270	1010	1018	1033	rBV	655871	1302398	18.49%	5.005%
8	10.915	1290	1298	1317	rBV	273787	589182	8.37%	2.264%
9	13.348	1704	1712	1725	rBV	2196439	3603771	51.17%	13.848%
10	14.728	1939	1947	1964	rVB2	504544	905939	12.86%	3.481%
11	16.215	2193	2200	2219	rBV	1755466	2899942	41.18%	11.143%
12	17.466	2406	2413	2430	rBV2	773186	1291064	18.33%	4.961%
13	20.063	2848	2855	2871	rBV	5309590	7042628	100.00%	27.062%
14	21.738	3133	3140	3154	rBV	960415	1706767	24.23%	6.559%
15	25.022	3687	3699	3714	rBV3	461020	1447840	20.56%	5.564%

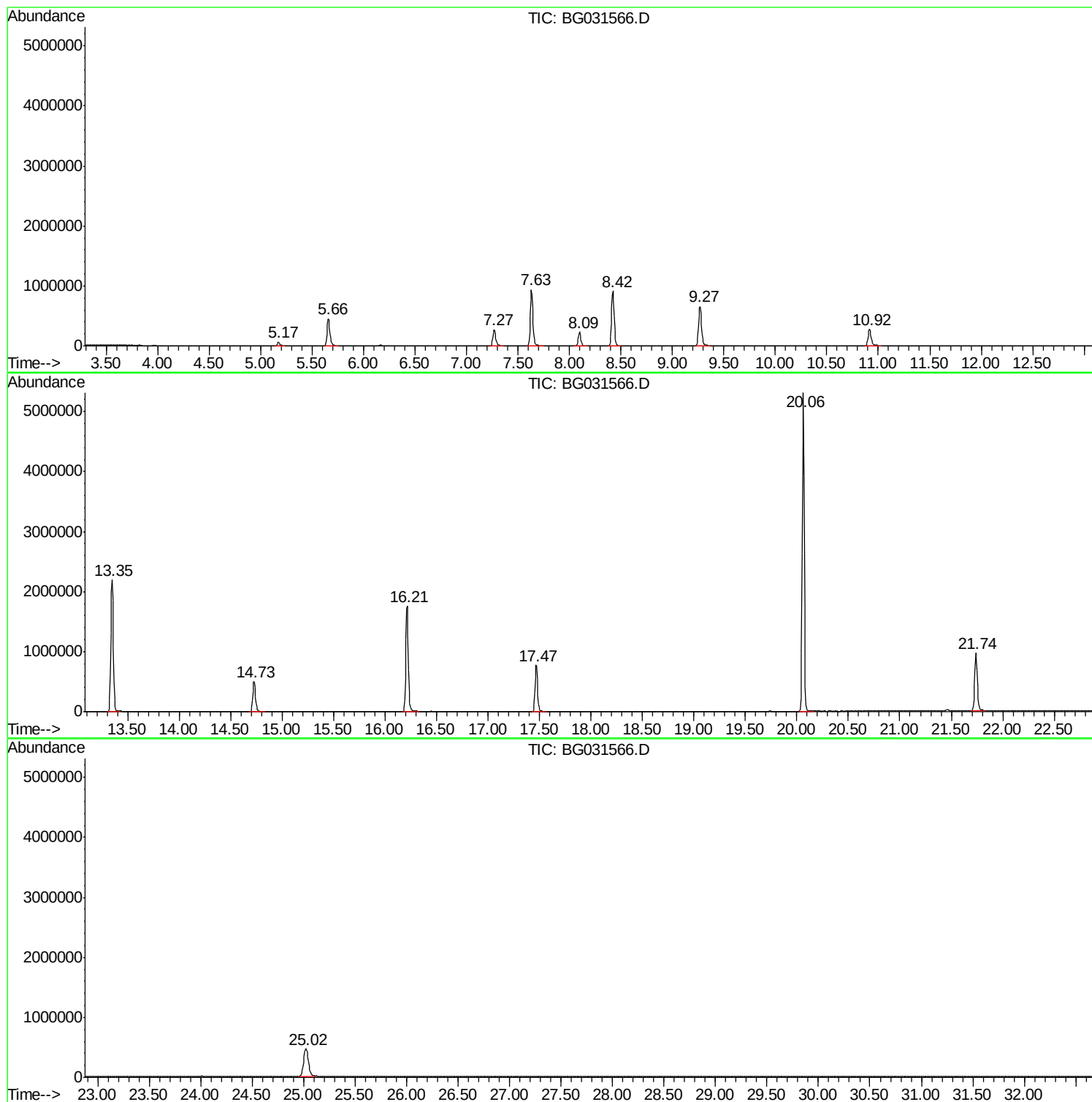
Sum of corrected areas: 26023621

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.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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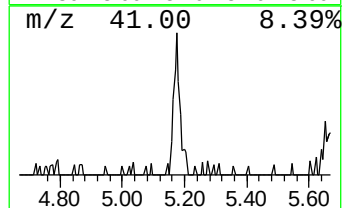
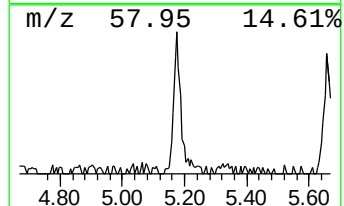
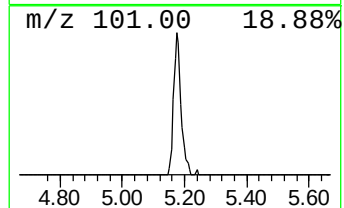
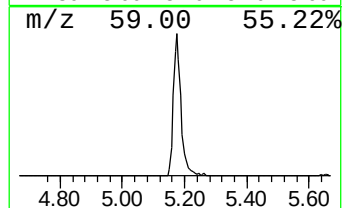
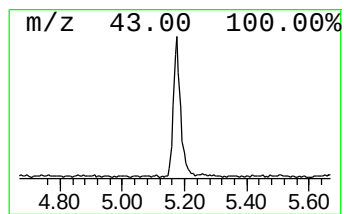
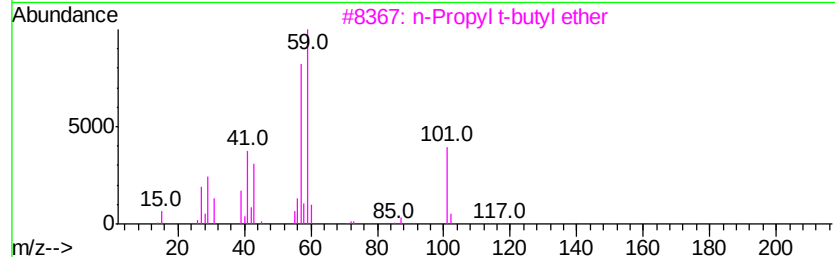
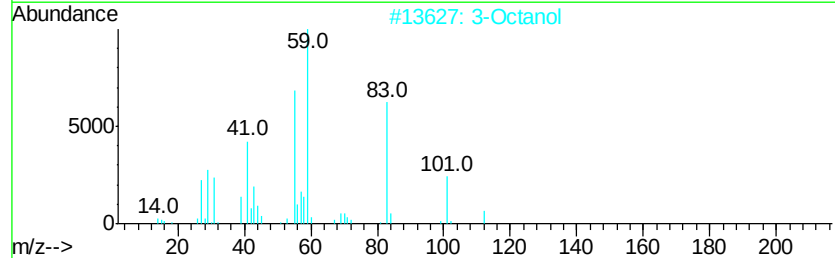
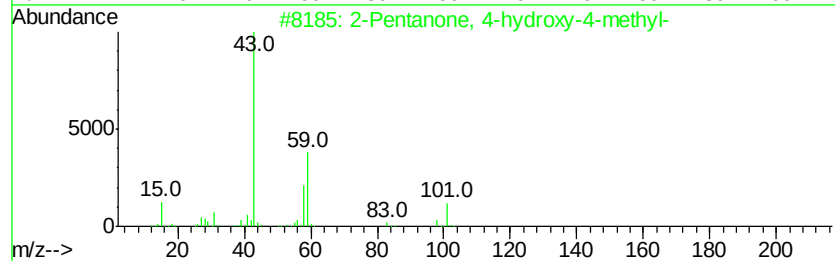
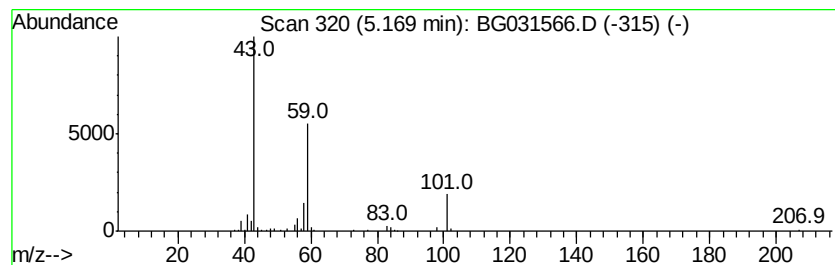
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.03 ng	104338	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Octanol	130	C8H18O	000589-98-0	38
3			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25



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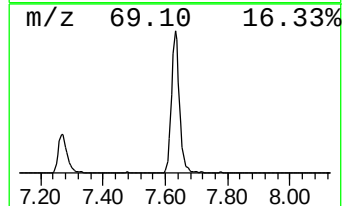
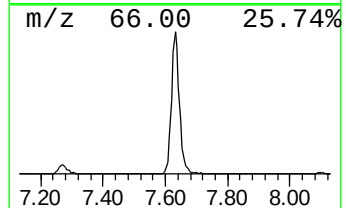
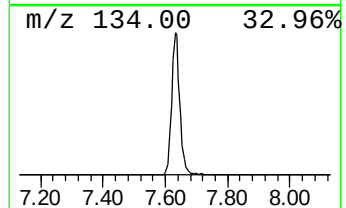
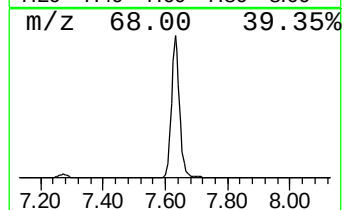
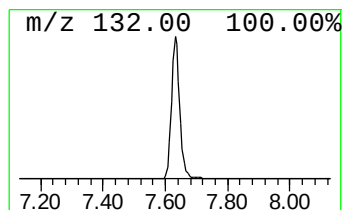
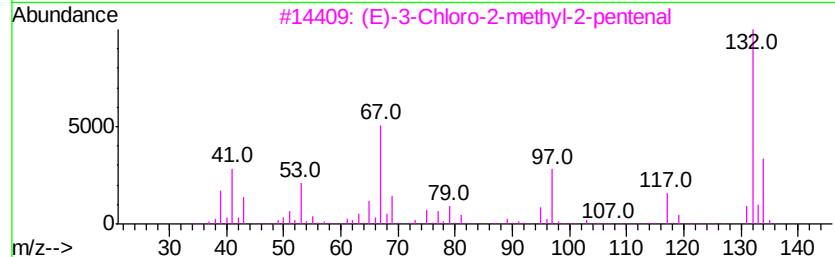
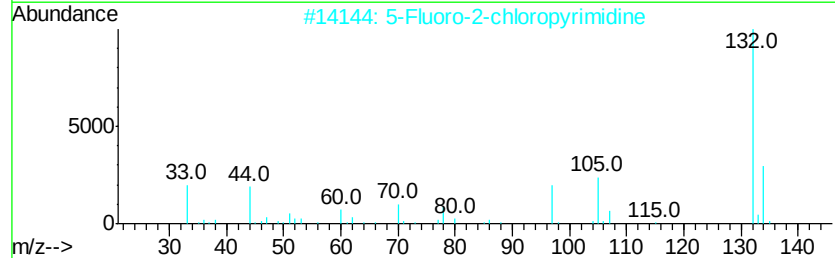
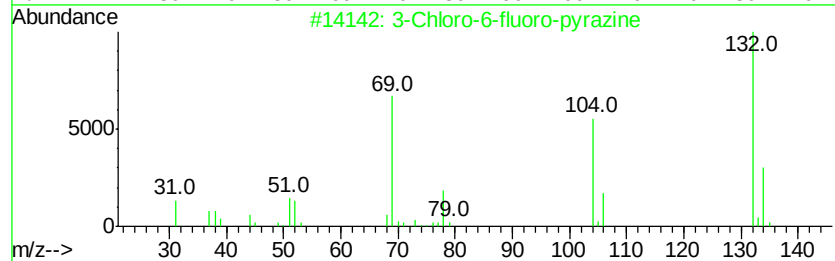
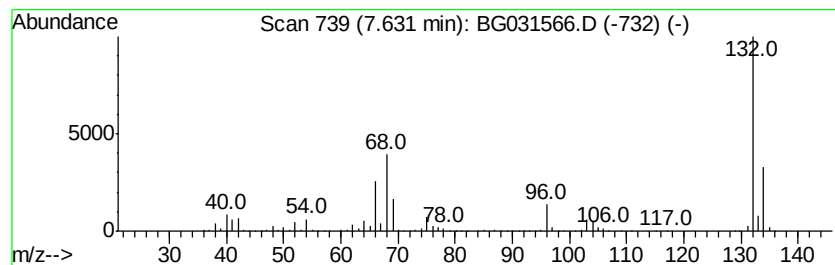
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	81.52 ng	1691170	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	5.0	ng	104338	1	8.09	414907	20.0
unknown7.63	7.63	81.5	ng	1691170	1	8.09	414907	20.0