

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031474.D
 Acq On : 12 Dec 2017 3:50
 Operator : SJ/JU
 Sample : I6762-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120417-JETT-E-D-S

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	5.185	318	323	331	rBV	46166	73995	1.30%	0.394%
2	5.667	399	405	422	rBV	243469	448980	7.86%	2.392%
3	7.277	669	679	690	rBV	155994	332920	5.83%	1.774%
4	7.641	732	741	759	rBV	479931	884268	15.48%	4.712%
5	8.111	814	821	829	rBV	159990	277393	4.86%	1.478%
6	8.434	868	876	886	rBV	692086	1256691	22.00%	6.696%
7	9.281	1012	1020	1033	rBV	463454	950243	16.64%	5.063%
8	10.932	1293	1301	1318	rBV	168258	396639	6.94%	2.113%
9	13.352	1706	1713	1737	rBV	1616072	2781574	48.70%	14.821%
10	14.733	1940	1948	1962	rVB2	374188	667256	11.68%	3.555%
11	16.219	2194	2201	2215	rBV	875262	1462818	25.61%	7.795%
12	17.477	2408	2415	2429	rBV2	580382	981519	17.18%	5.230%
13	20.068	2849	2856	2869	rBV	4216543	5711764	100.00%	30.435%
14	21.355	3070	3075	3083	rBV9	22296	60754	1.06%	0.324%
15	21.742	3134	3141	3156	rBV2	750299	1293866	22.65%	6.894%
16	25.021	3688	3699	3715	rBV3	397506	1186613	20.77%	6.323%

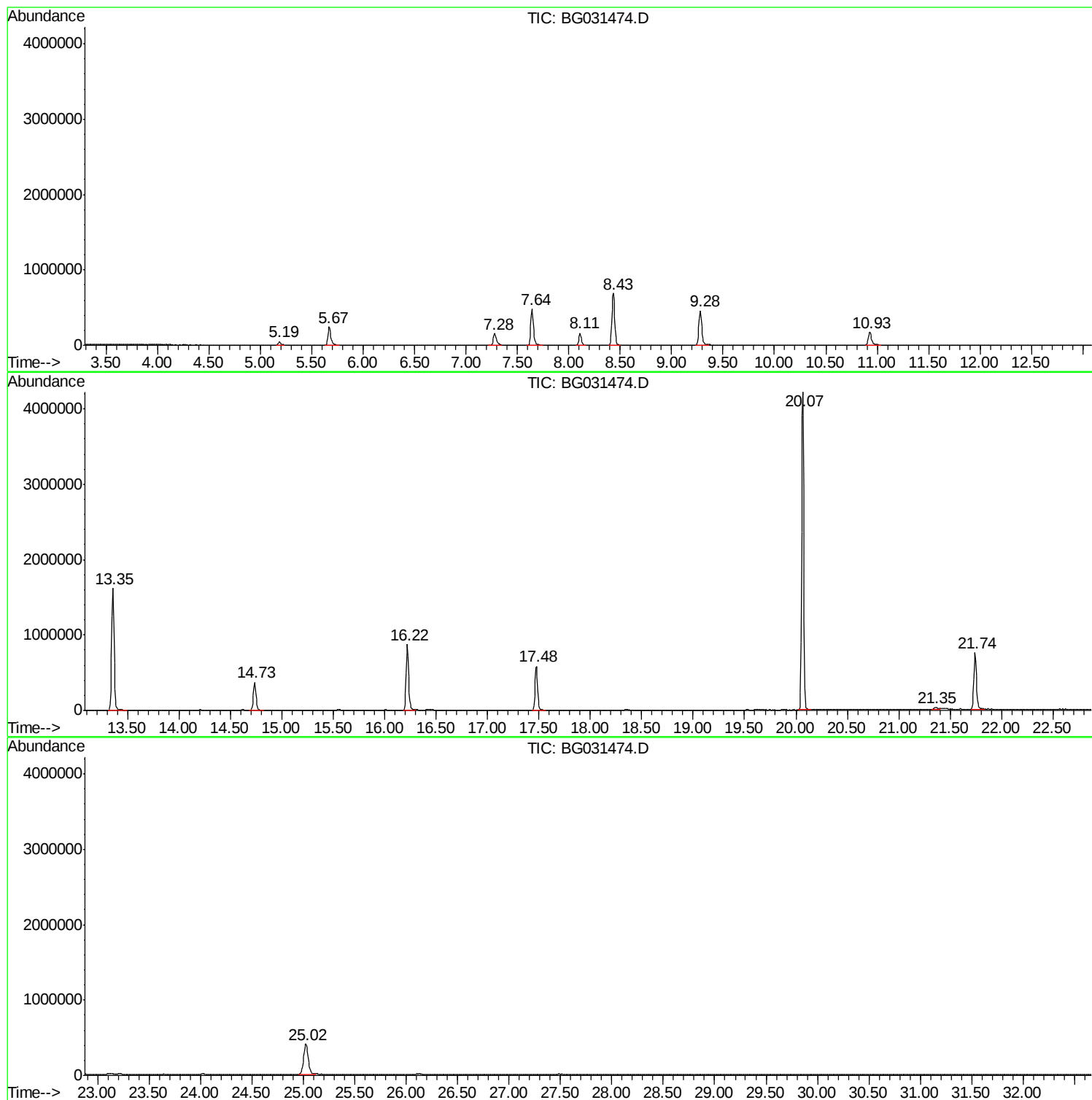
Sum of corrected areas: 18767293

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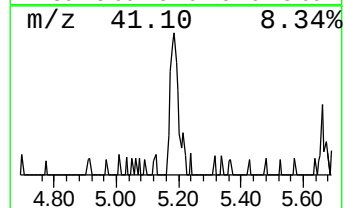
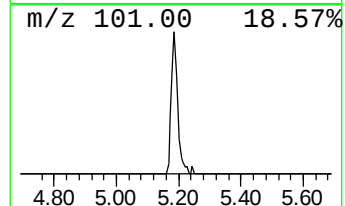
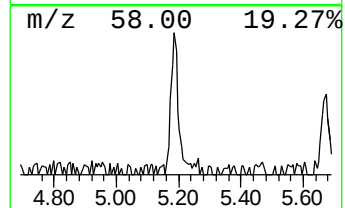
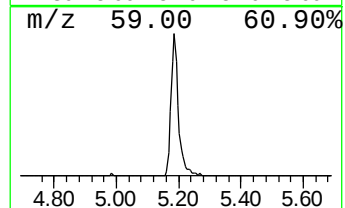
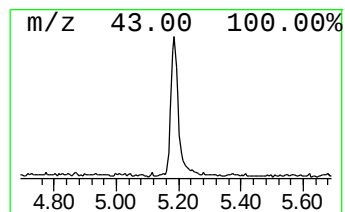
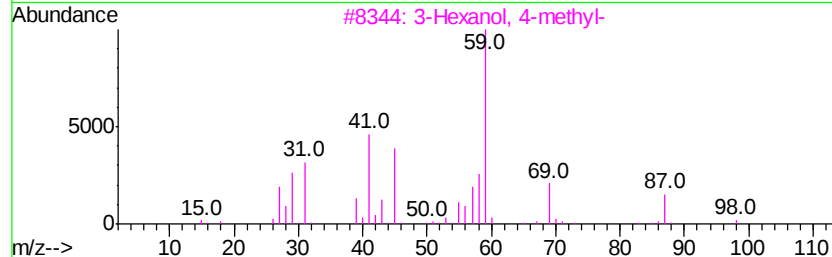
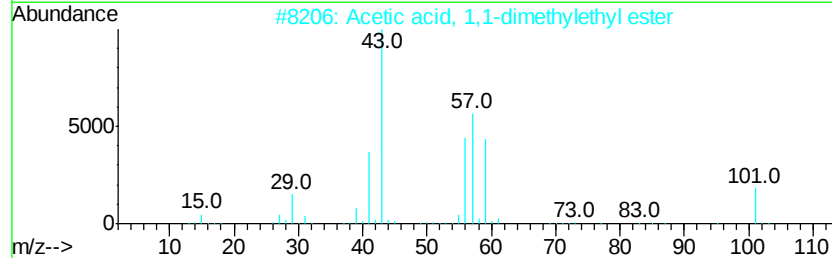
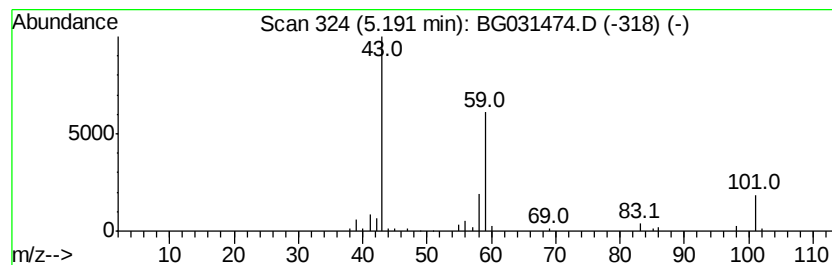
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	5.34 ng	73995	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
4			1,2-Butanediol	90	C4H10O2	000584-03-2	17
5			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	17



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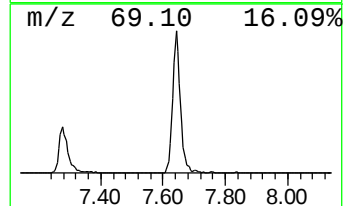
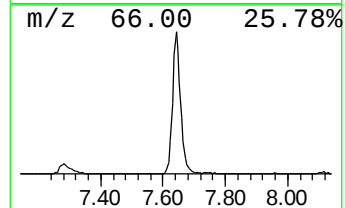
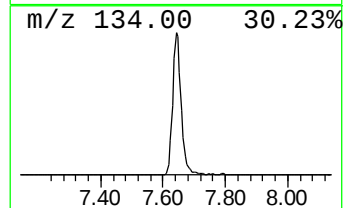
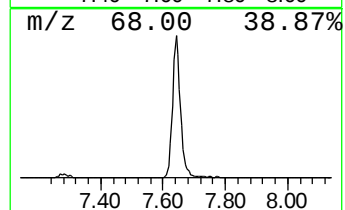
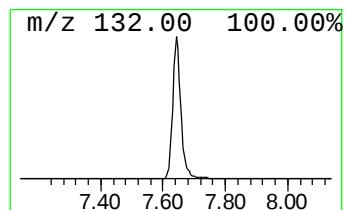
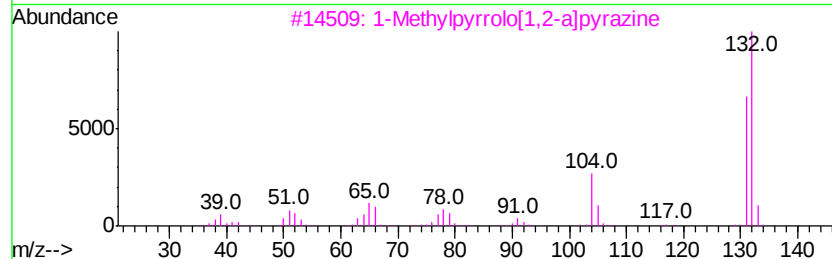
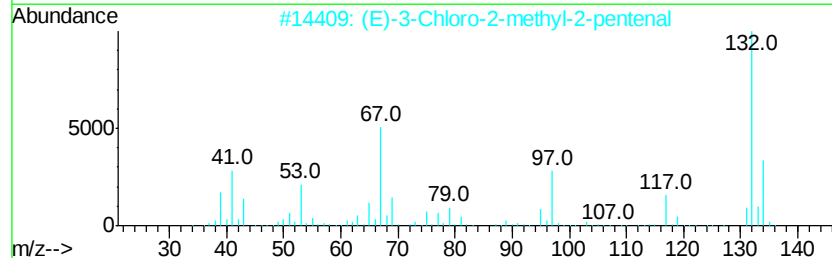
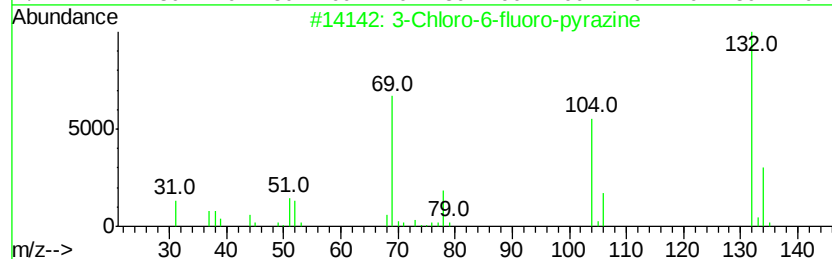
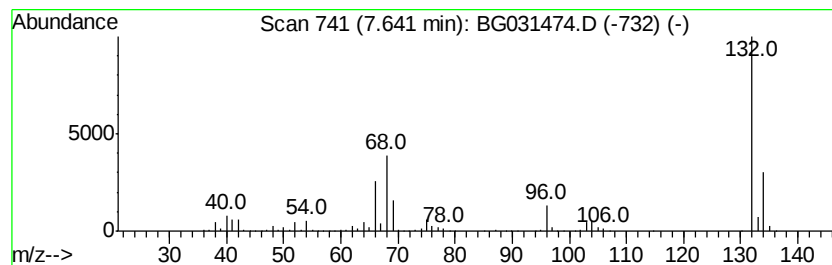
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	63.76 ng	884268	1,4-Dichlorobenzene-d4	8.11

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	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	5.3	ng	73995	1	8.11	277393	20.0
unknown7.64	7.64	63.8	ng	884268	1	8.11	277393	20.0