

Data Path : Z:\HPCHEM1\BNA G\DATA\BG082316\
 Data File : BG023860.D
 Acq On : 23 Aug 2016 19:26
 Operator : UM/SJ
 Sample : H4556-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EB-01

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-BG080116.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	5.152	389	394	405	rVB	169433	273033	3.33%	0.748%
2	5.628	469	475	490	rBV	826999	1437051	17.53%	3.939%
3	7.244	744	750	764	rBV	533847	965587	11.78%	2.647%
4	7.619	805	814	824	rBV	1545035	2798128	34.13%	7.670%
5	8.089	885	894	902	rVB2	377234	648754	7.91%	1.778%
6	8.413	938	949	958	rBV	1445163	2572562	31.38%	7.052%
7	9.270	1086	1095	1106	rBV	1265006	2343398	28.58%	6.424%
8	10.927	1367	1377	1386	rBV	479351	909677	11.10%	2.494%
9	13.376	1785	1794	1806	rBV	3396732	5268652	64.27%	14.443%
10	14.763	2023	2030	2043	rBV2	801474	1300085	15.86%	3.564%
11	16.261	2278	2285	2300	rBV	2640902	4418864	53.90%	12.113%
12	17.524	2494	2500	2515	rVB	1056630	1661189	20.26%	4.554%
13	20.132	2938	2944	2951	rBV	6033346	8198269	100.00%	22.473%
14	21.841	3228	3235	3243	rVB	1082169	1871693	22.83%	5.131%
15	25.213	3797	3809	3822	rBV4	564283	1812851	22.11%	4.969%

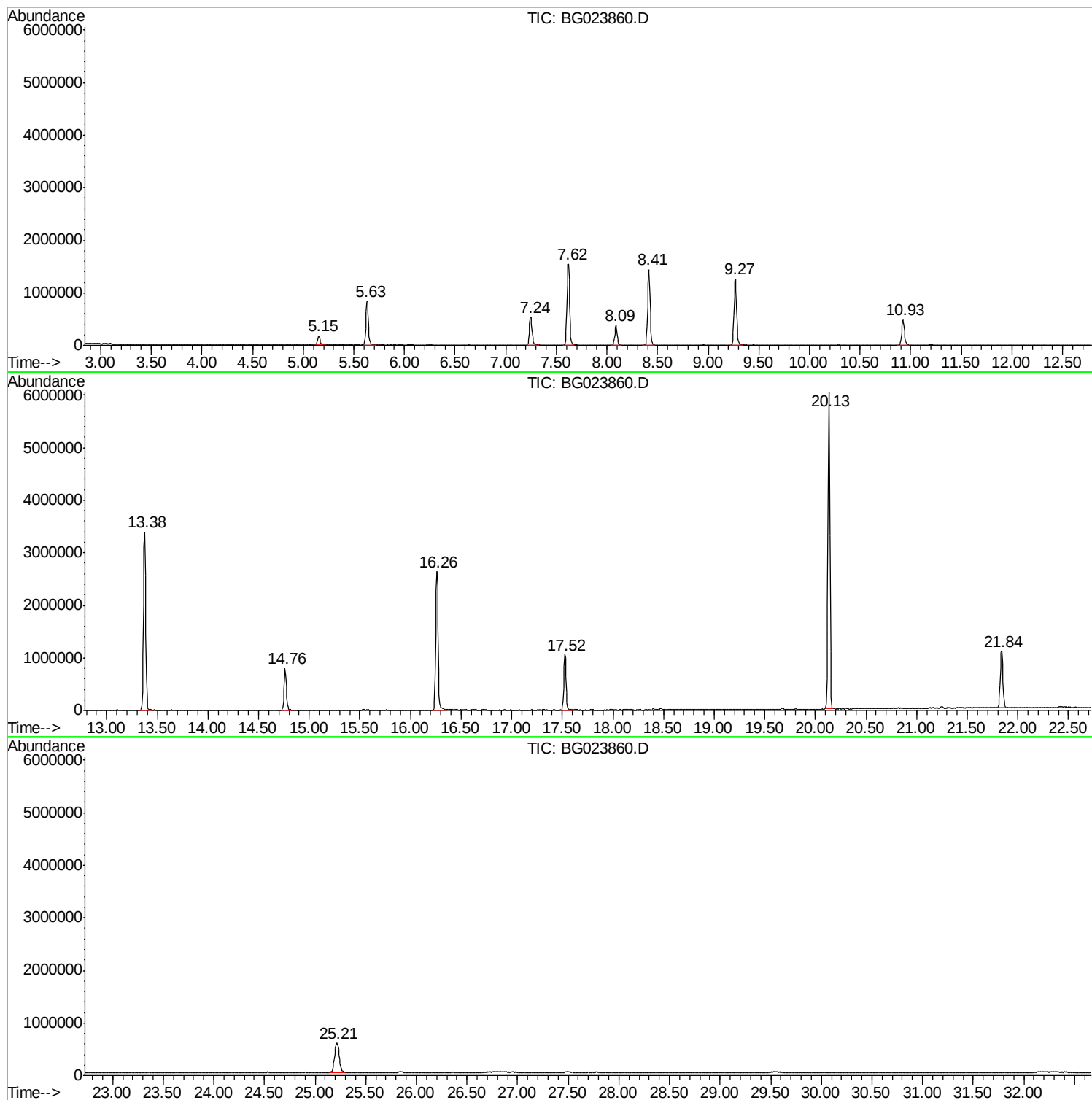
Sum of corrected areas: 36479793

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Quant Method : Z:\HPCHEM1\BNA G\Methods\8270-BG080116.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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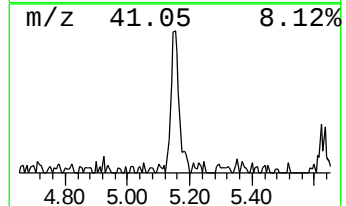
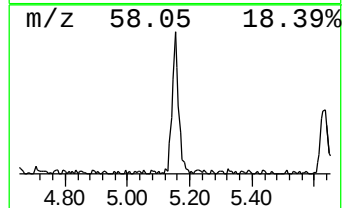
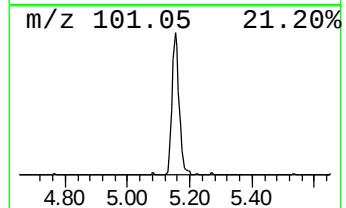
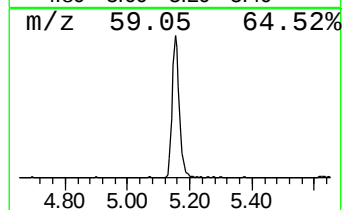
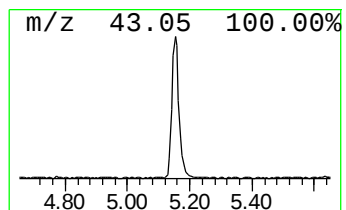
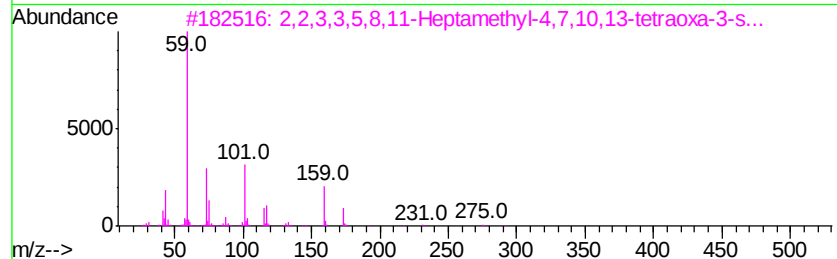
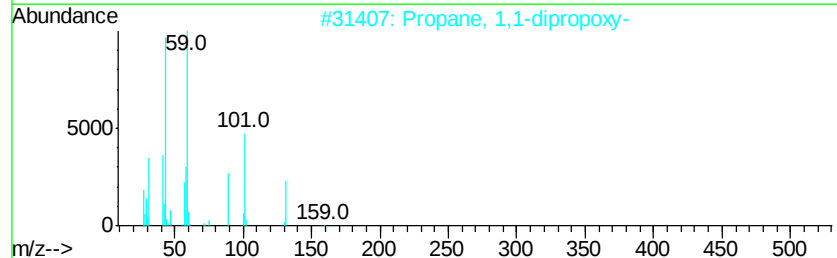
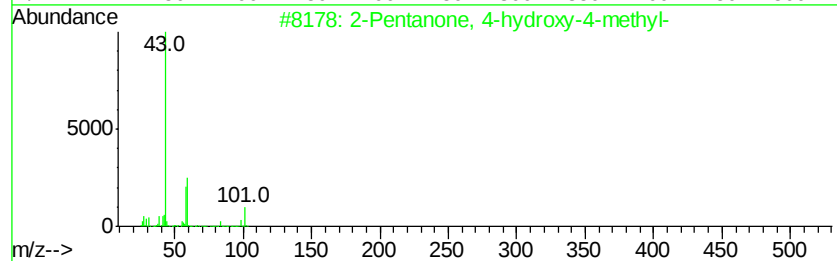
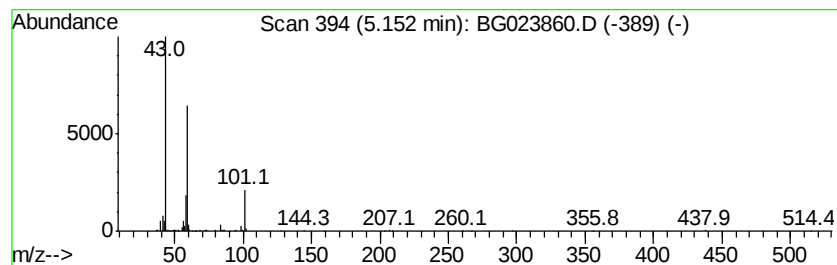
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	8.42 ng	273033	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	59
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
3			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	28
4			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	23
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23



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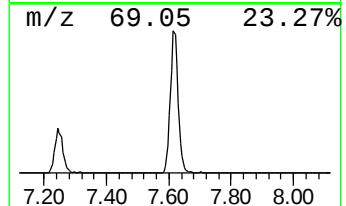
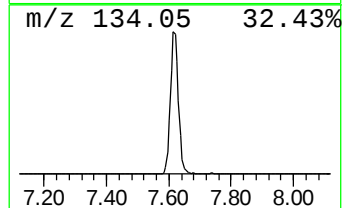
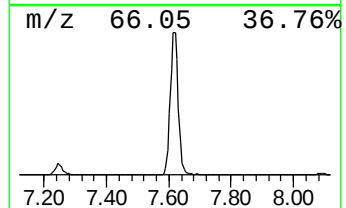
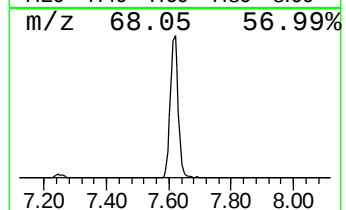
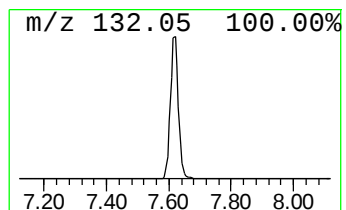
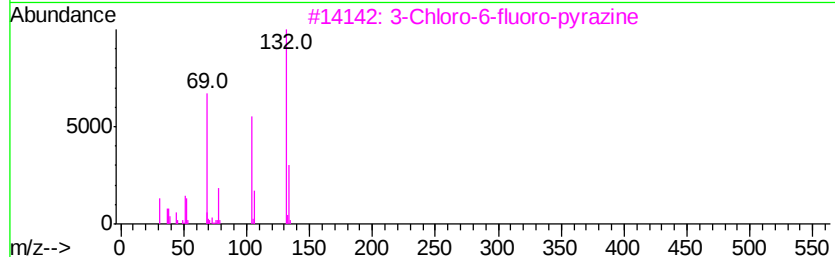
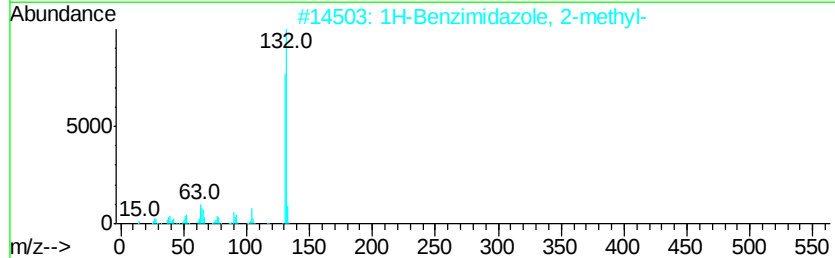
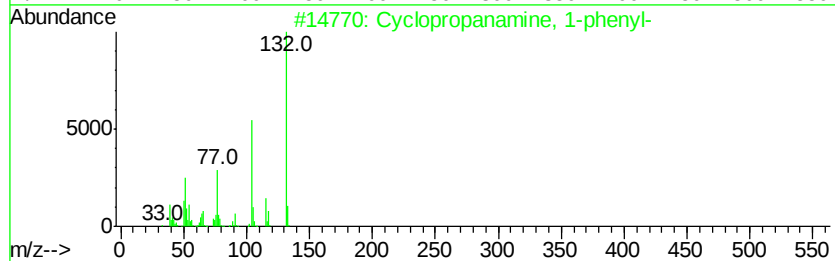
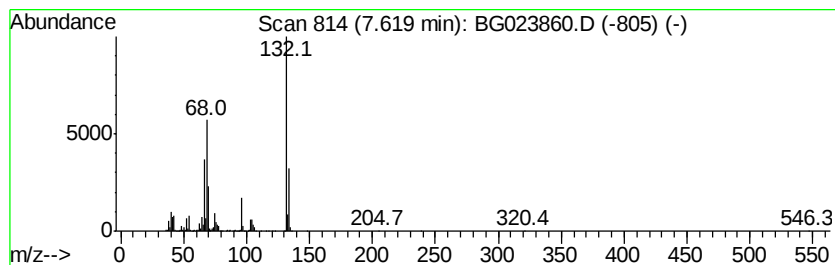
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	86.26 ng	2798130	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	8.4	ng	273033	1	8.09	648754	20.0
unknown7.62	7.62	86.3	ng	2798130	1	8.09	648754	20.0