

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061818\
 Data File : BG035222.D
 Acq On : 18 Jun 2018 17:13
 Operator : SJ/JU
 Sample : J3464-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-42-COMP

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-BG060718.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	316	321	330	rBV2	111130	177334	2.65%	0.848%
2	5.638	393	400	411	rBV	746517	1221751	18.24%	5.840%
3	7.218	661	669	679	rBV	702368	1207910	18.03%	5.774%
4	7.594	724	733	742	rBV	789374	1363401	20.35%	6.517%
5	8.064	806	813	821	rBB2	138593	239691	3.58%	1.146%
6	8.387	860	868	875	rBV	591859	1057710	15.79%	5.056%
7	9.221	1002	1010	1022	rVB	514596	910779	13.60%	4.354%
8	10.866	1283	1290	1295	rBV	169053	307662	4.59%	1.471%
9	10.919	1295	1299	1305	rVB2	50612	82618	1.23%	0.395%
10	13.310	1698	1706	1716	rBV	1300528	2081557	31.08%	9.950%
11	14.684	1932	1940	1948	rBV2	265792	448171	6.69%	2.142%
12	16.165	2186	2192	2209	rBV	1360384	2111250	31.52%	10.092%
13	17.428	2399	2407	2413	rBV	473739	744139	11.11%	3.557%
14	20.036	2844	2851	2856	rBV	4935674	6698459	100.00%	32.019%
15	21.692	3126	3133	3142	rBV2	773978	1264653	18.88%	6.045%
16	24.917	3671	3682	3697	rVB	350123	1002927	14.97%	4.794%

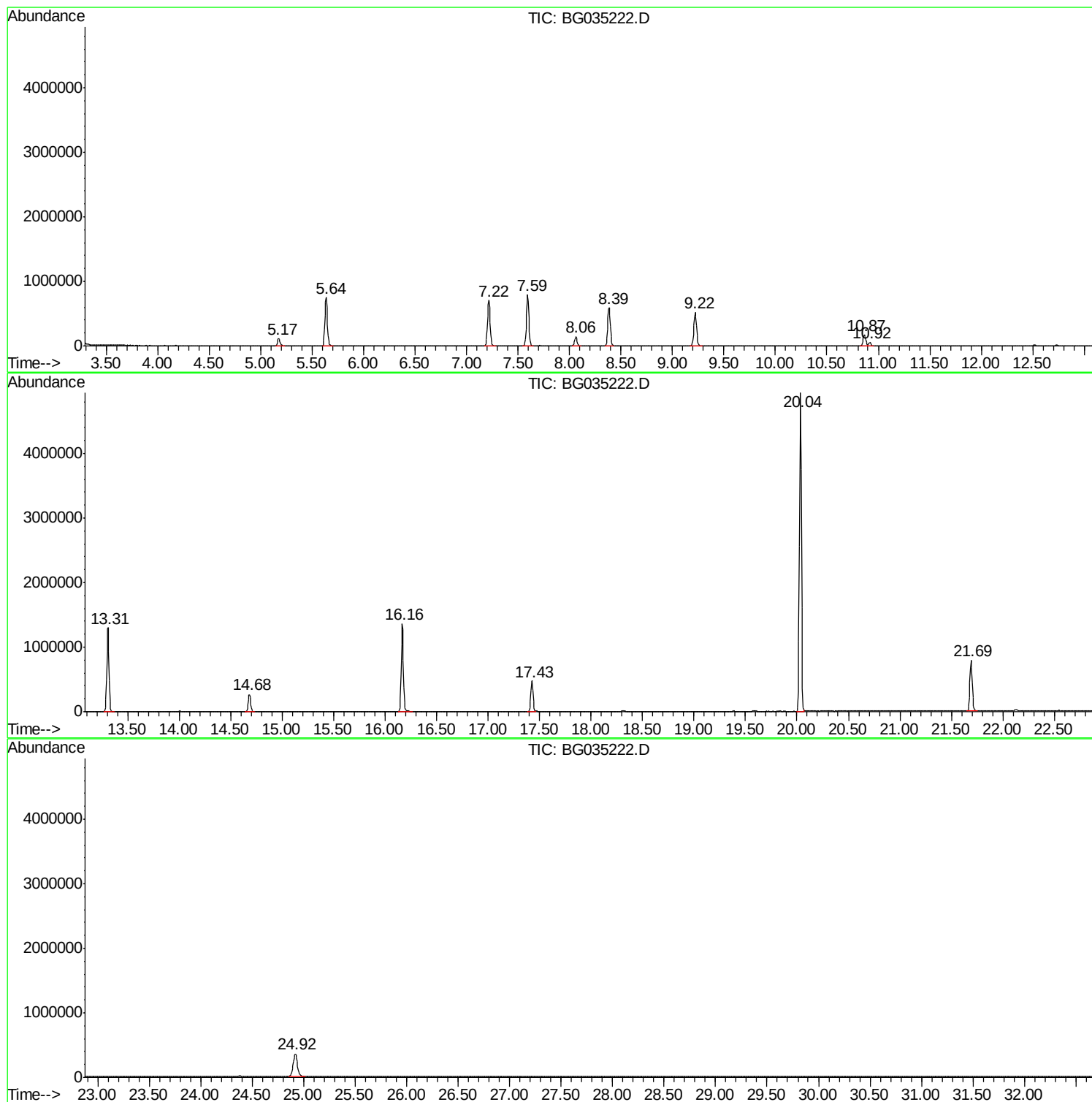
Sum of corrected areas: 20920012

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.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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ClientSampled :
SB-42-COMP

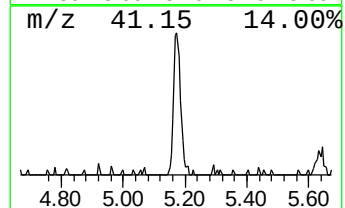
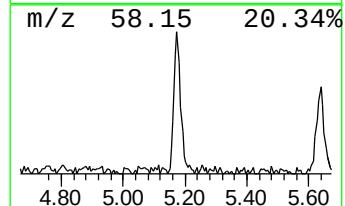
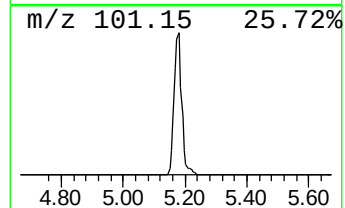
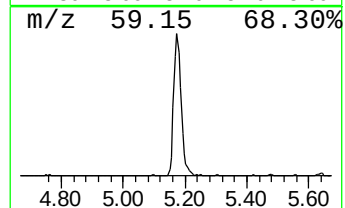
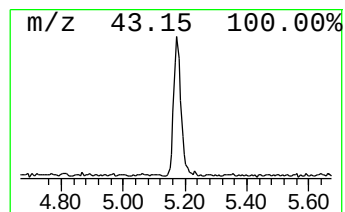
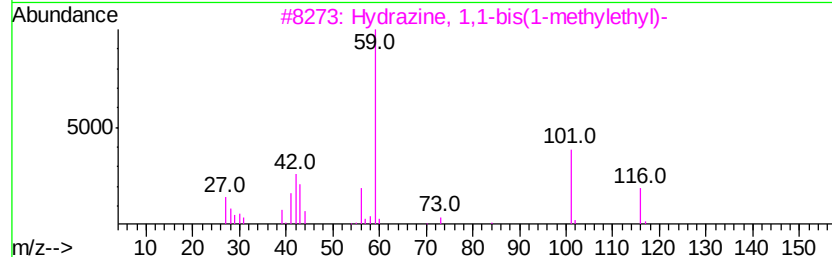
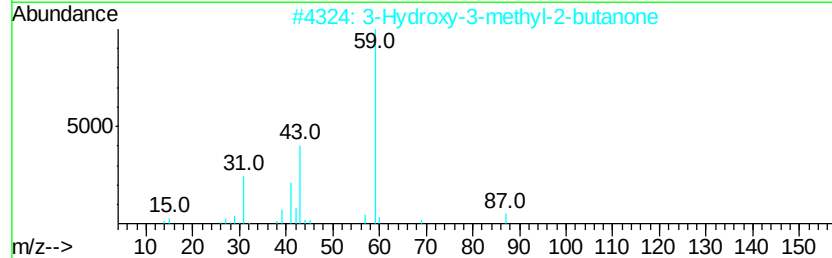
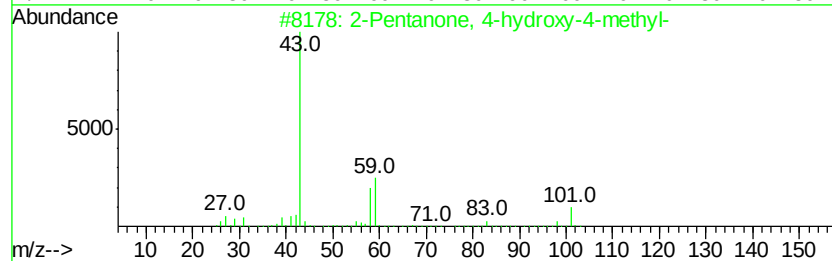
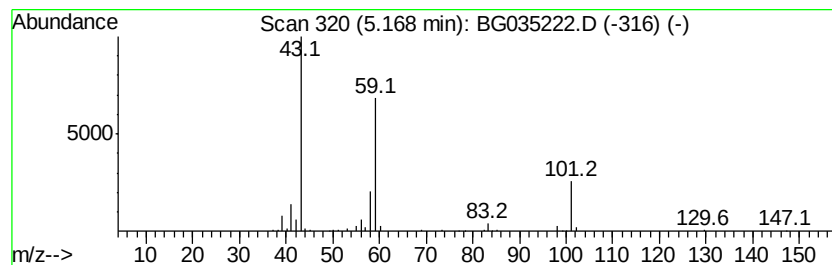
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	14.80 ng	177334	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	27
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
4			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	25
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	12



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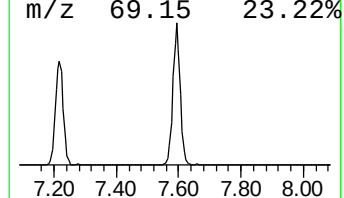
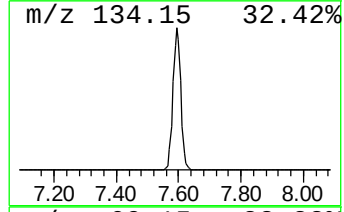
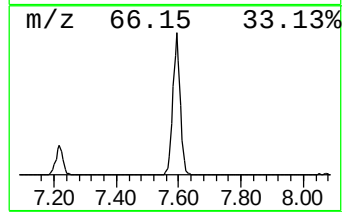
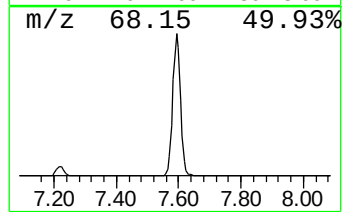
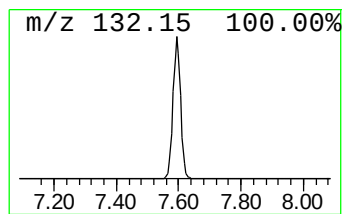
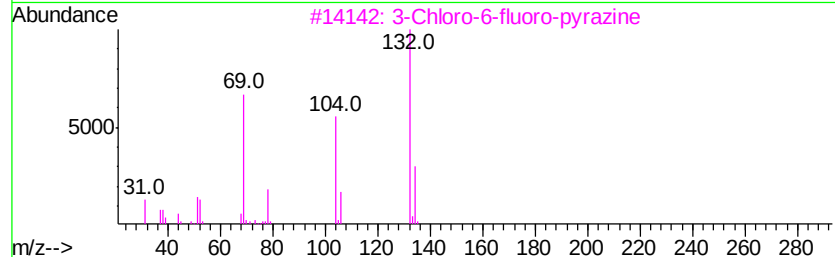
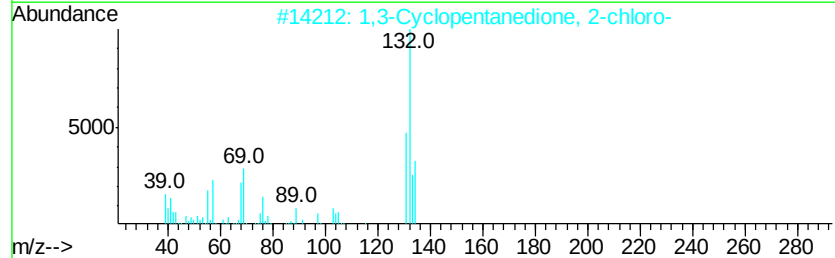
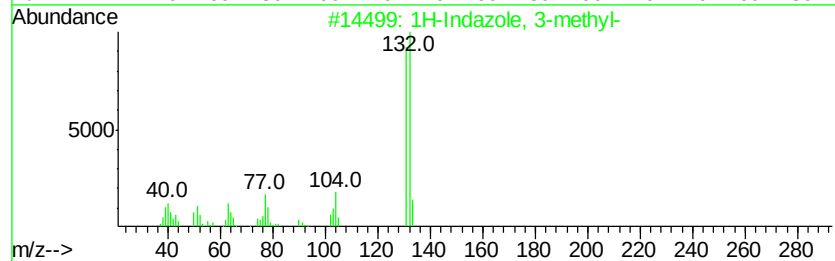
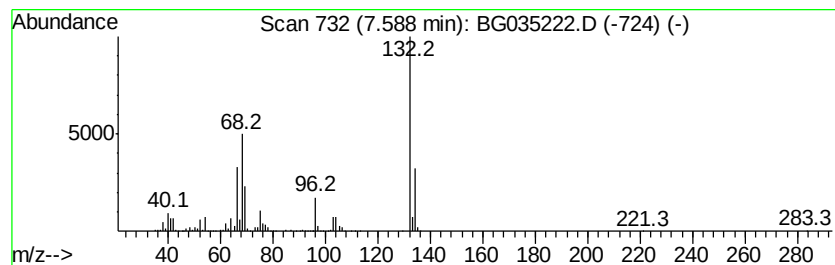
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	113.76 ng	1363400	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indazole, 3-methyl-	132	C8H8N2	1000316-00-2	14
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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	14.8	ng	177334	1	8.06	239691	20.0
unknown7.59	7.59	113.8	ng	1363400	1	8.06	239691	20.0