

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062117\
 Data File : BG027569.D
 Acq On : 21 Jun 2017 20:47
 Operator : SJ/JU
 Sample : PB99877BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99877BL

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-BG062117.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	5.640	292	298	307	rVB	61858	103594	3.53%	0.854%
2	6.087	366	374	385	rBV	563661	974117	33.23%	8.032%
3	7.644	631	639	654	rBV	560416	998851	34.08%	8.236%
4	8.038	698	706	719	rVB	534765	973127	33.20%	8.024%
5	8.508	778	786	794	rVB	87932	158484	5.41%	1.307%
6	8.831	833	841	850	rBV	375158	681439	23.25%	5.619%
7	9.671	975	984	992	rBV	307409	581201	19.83%	4.792%
8	11.322	1258	1265	1276	rVB	119481	228015	7.78%	1.880%
9	13.713	1665	1672	1686	rBV	848321	1318474	44.98%	10.872%
10	15.088	1900	1906	1914	rBV2	181496	304047	10.37%	2.507%
11	16.575	2152	2159	2171	rBV	804437	1248456	42.59%	10.295%
12	17.838	2366	2374	2382	rBV2	293048	466837	15.93%	3.849%
13	20.417	2806	2813	2819	rBV	2243813	2931323	100.00%	24.171%
14	22.209	3110	3118	3127	rBV	321183	596448	20.35%	4.918%
15	25.846	3724	3737	3750	rVB2	175325	562976	19.21%	4.642%

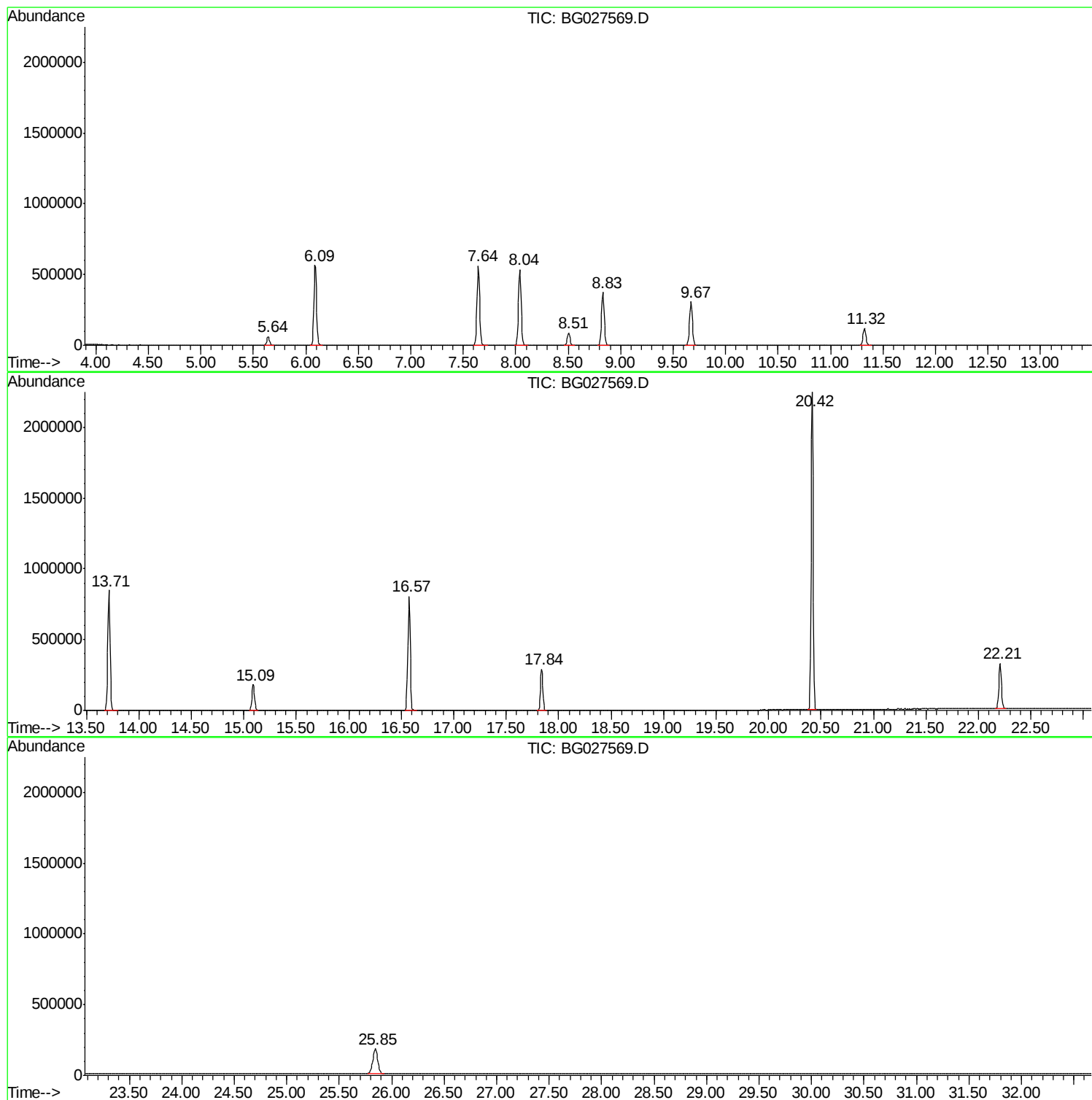
Sum of corrected areas: 12127389

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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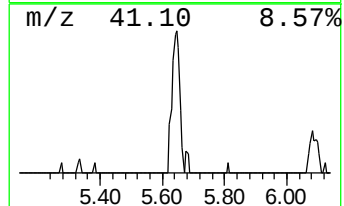
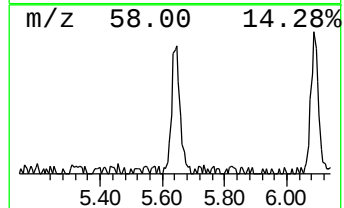
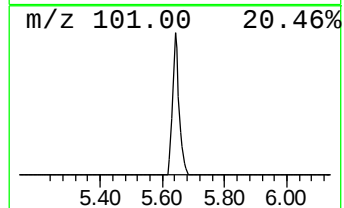
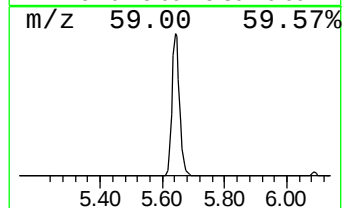
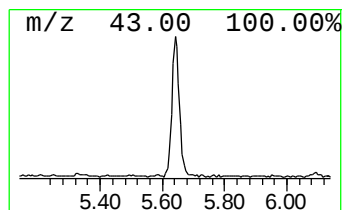
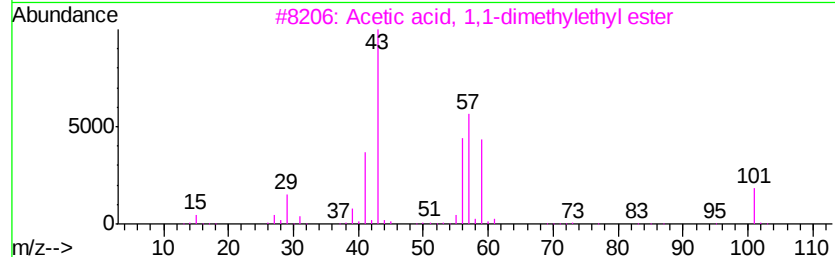
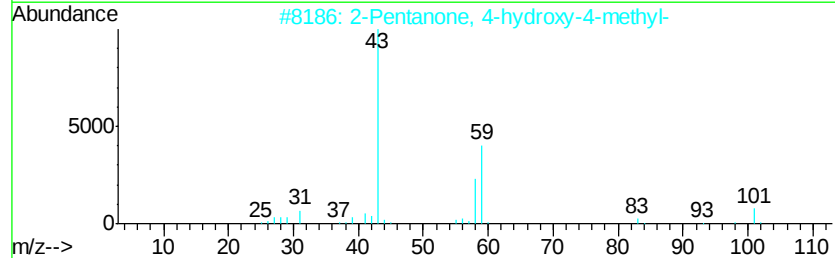
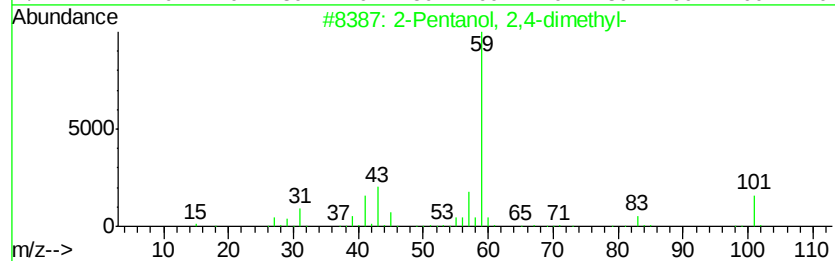
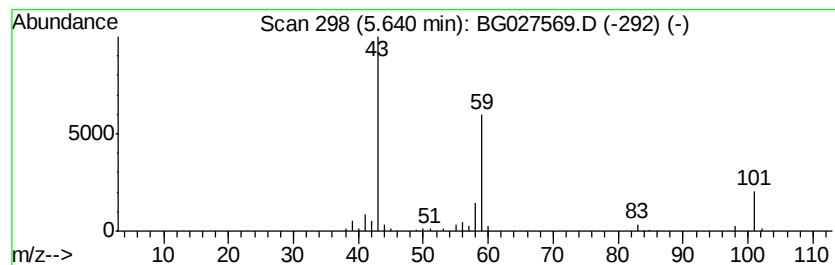
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanol, 2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	13.07 ng	103594	1,4-Dichlorobenzene-d4	8.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	53
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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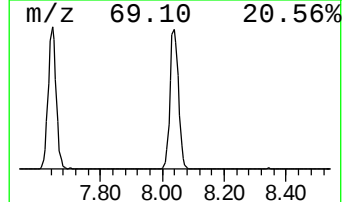
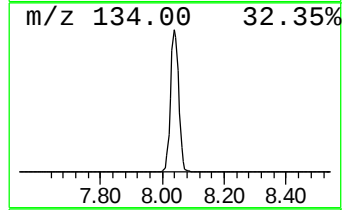
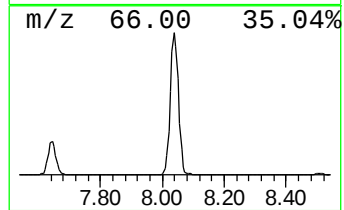
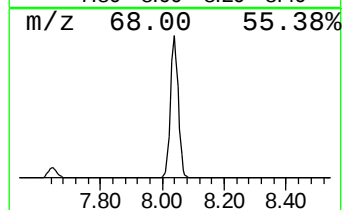
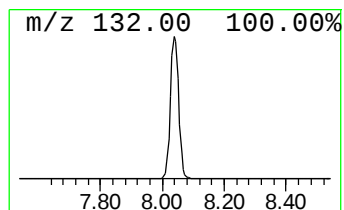
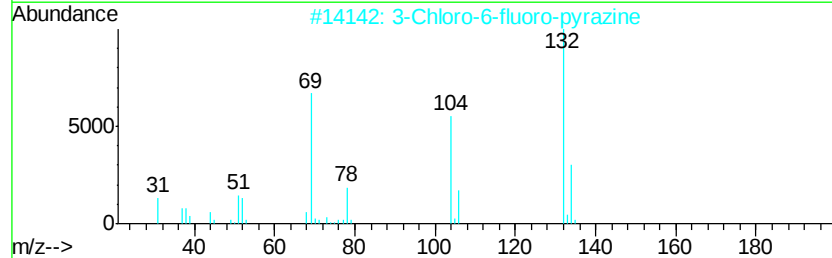
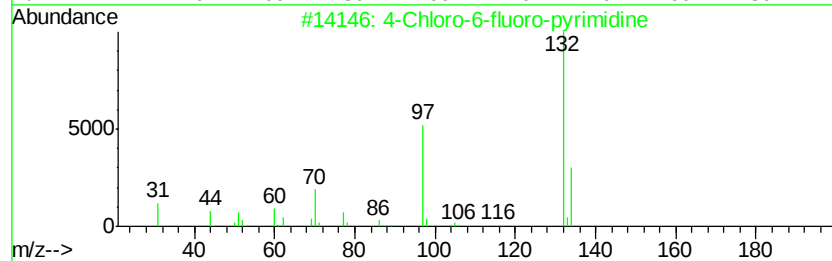
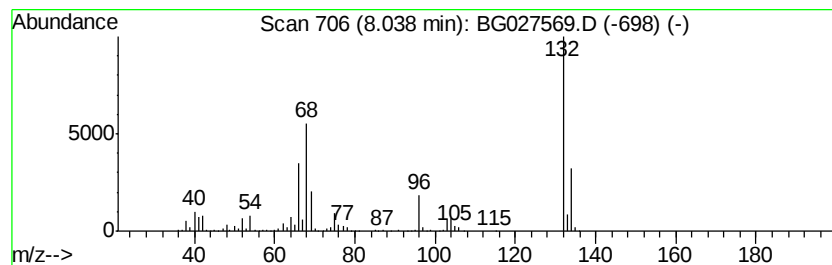
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Peak Number 2 unknown8.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	122.80 ng	973127	1,4-Dichlorobenzene-d4	8.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanol, 2,4-d...	5.64	13.1	ng	103594	1	8.51	158484	20.0
unknown8.04	8.04	122.8	ng	973127	1	8.51	158484	20.0