

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110718\
 Data File : BG037875.D
 Acq On : 7 Nov 2018 21:57
 Operator : JU/SJ
 Sample : PB114524BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114524BL

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-BG101818.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	5.171	315	320	330	rVB	92472	151308	4.80%	0.894%
2	5.635	392	399	414	rVB	788257	1310904	41.58%	7.743%
3	7.239	663	672	685	rBV	852521	1506265	47.77%	8.896%
4	7.621	729	737	747	rBV	790947	1412907	44.81%	8.345%
5	8.091	809	817	825	rBV	156551	280328	8.89%	1.656%
6	8.414	864	872	884	rVB	585134	1057982	33.56%	6.249%
7	9.272	1008	1018	1027	rBV	491123	927091	29.40%	5.476%
8	10.911	1290	1297	1307	rVB	232218	437887	13.89%	2.586%
9	13.350	1704	1712	1723	rBV	1468064	2355276	74.70%	13.911%
10	14.724	1939	1946	1955	rBV2	386185	656063	20.81%	3.875%
11	16.217	2192	2200	2219	rBV	947838	1542523	48.92%	9.111%
12	17.474	2405	2414	2426	rBV	498156	769181	24.40%	4.543%
13	20.071	2850	2856	2873	rBV	2236074	3152967	100.00%	18.622%
14	21.757	3136	3143	3158	rBV	422335	773287	24.53%	4.567%
15	25.089	3698	3710	3726	rBV3	163119	597107	18.94%	3.527%

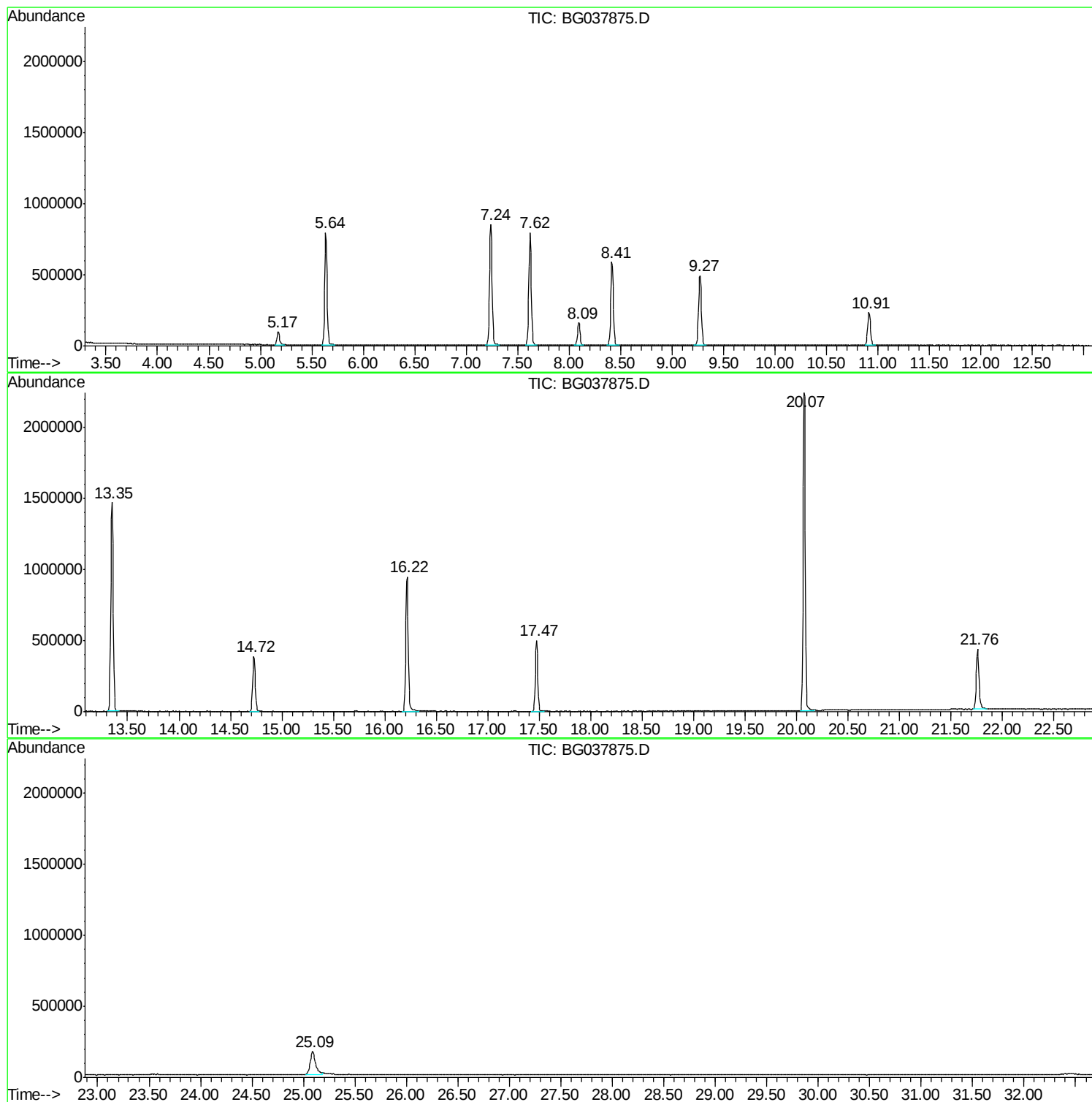
Sum of corrected areas: 16931076

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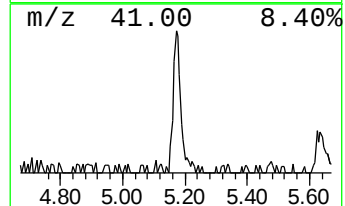
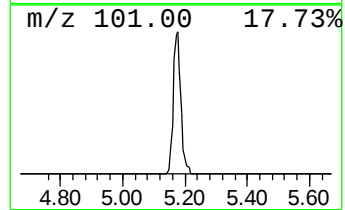
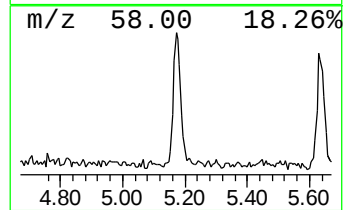
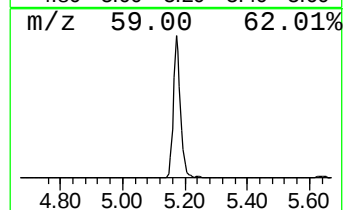
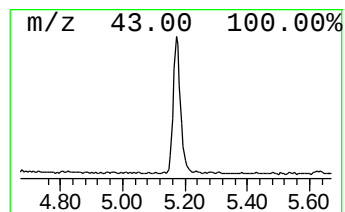
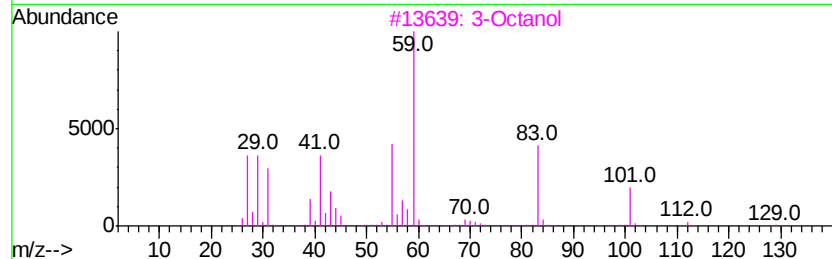
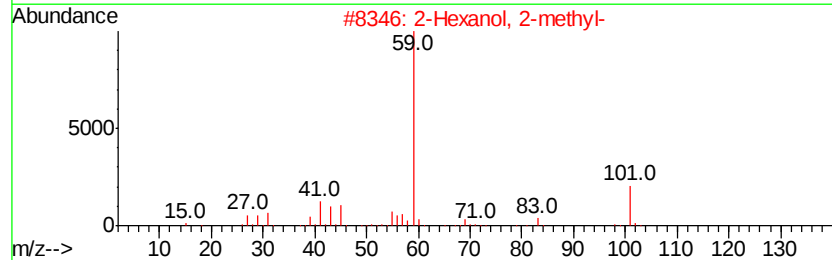
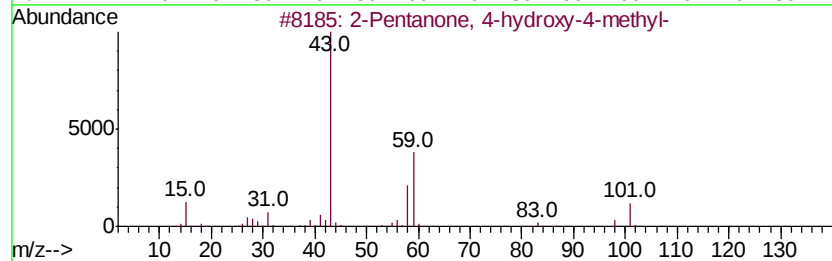
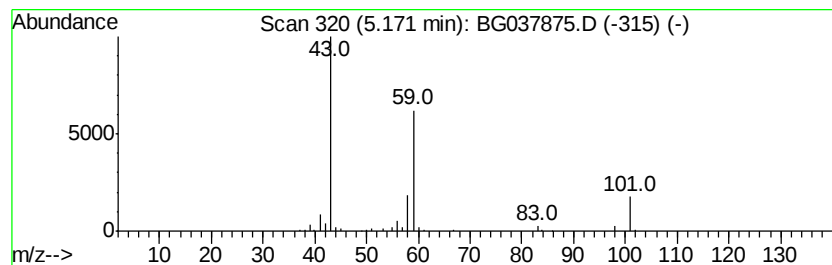
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	10.80 ng	151308	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		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3		3-Octanol	130	C8H18O	000589-98-0	33
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25



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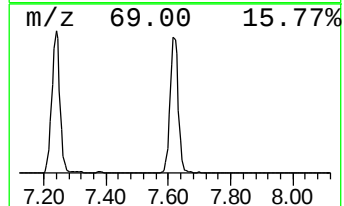
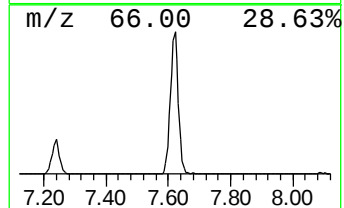
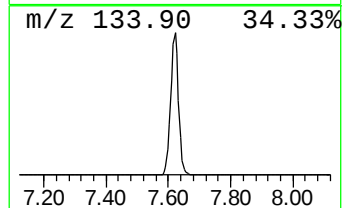
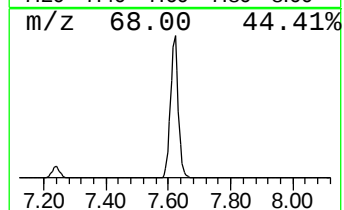
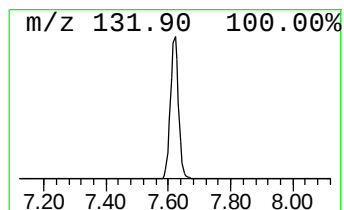
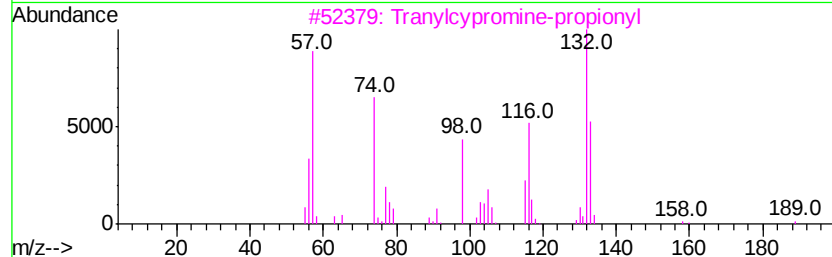
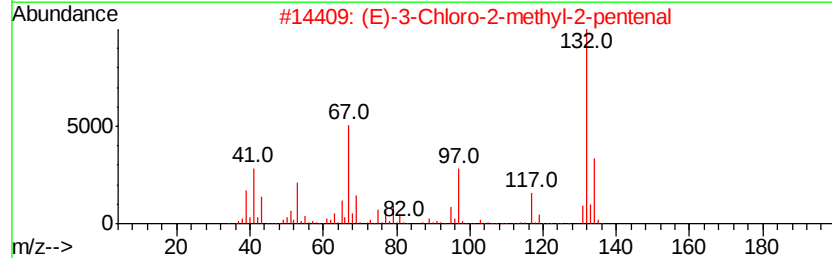
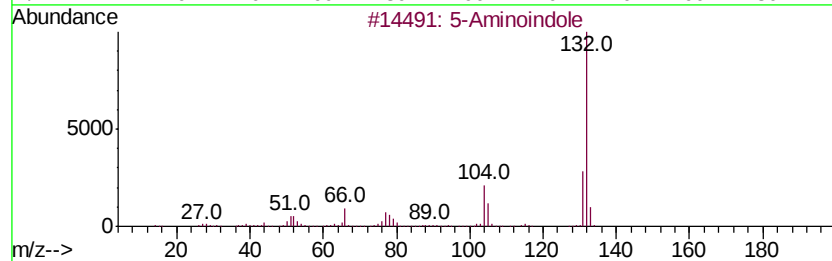
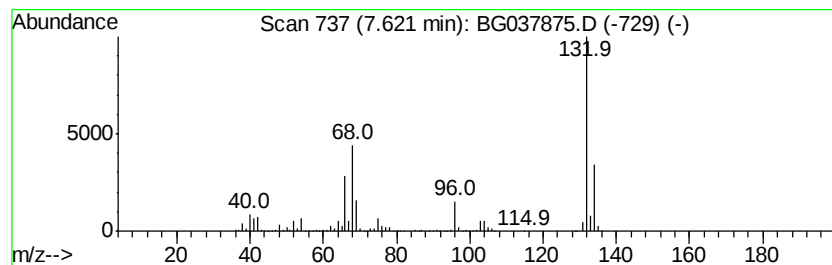
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	100.80 ng	1412910	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	10.8	ng	151308	1	8.09	280328	20.0
unknown7.62	7.62	100.8	ng	1412910	1	8.09	280328	20.0