

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103118\
 Data File : BG037684.D
 Acq On : 31 Oct 2018 15:00
 Operator : JU/SJ
 Sample : PB114332BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114332BL

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	5.180	317	322	328	rBV	28988	46025	1.64%	0.295%
2	5.644	393	401	411	rBV	644960	1086386	38.68%	6.964%
3	7.242	665	673	686	rBV	716553	1251658	44.57%	8.024%
4	7.624	729	738	748	rBV	650079	1152943	41.05%	7.391%
5	8.100	812	819	827	rVB	181473	316919	11.28%	2.032%
6	8.423	866	874	882	rBV	471568	831801	29.62%	5.332%
7	9.275	1012	1019	1034	rBV	399652	736238	26.21%	4.720%
8	10.920	1291	1299	1309	rBV	270592	497942	17.73%	3.192%
9	13.353	1706	1713	1723	rBV	1178878	1940132	69.08%	12.437%
10	14.733	1941	1948	1958	rBV2	487402	801946	28.55%	5.141%
11	16.220	2194	2201	2212	rBV	771253	1256841	44.75%	8.057%
12	17.483	2407	2416	2426	rBV	633492	1001537	35.66%	6.420%
13	20.080	2852	2858	2871	rBV	2092984	2808475	100.00%	18.004%
14	21.766	3137	3145	3158	rBV2	557547	1060276	37.75%	6.797%
15	25.098	3701	3712	3733	rBV2	207523	810161	28.85%	5.194%

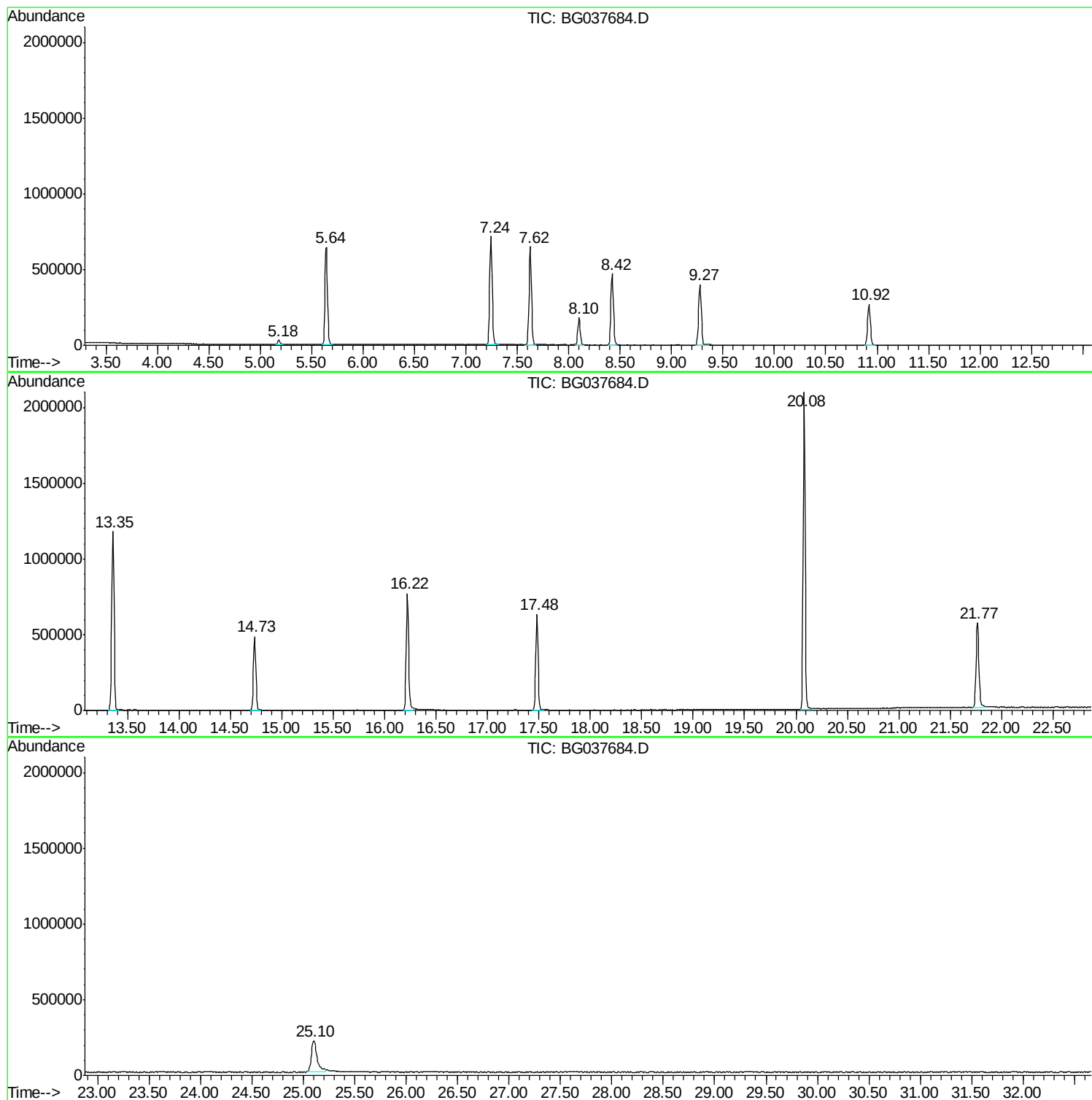
Sum of corrected areas: 15599280

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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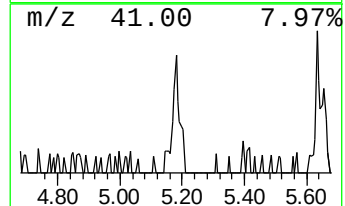
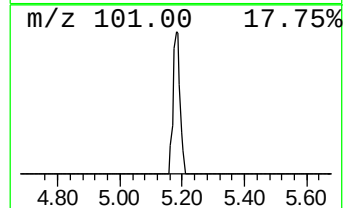
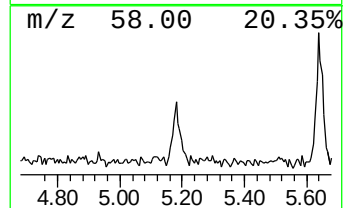
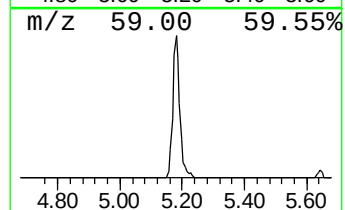
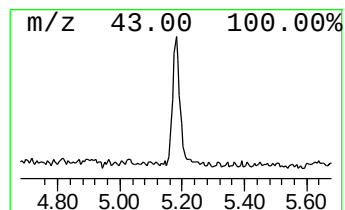
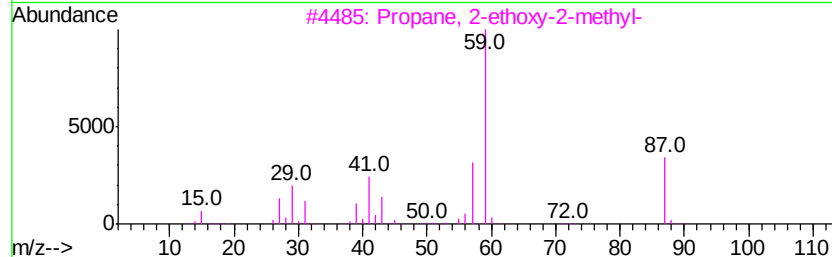
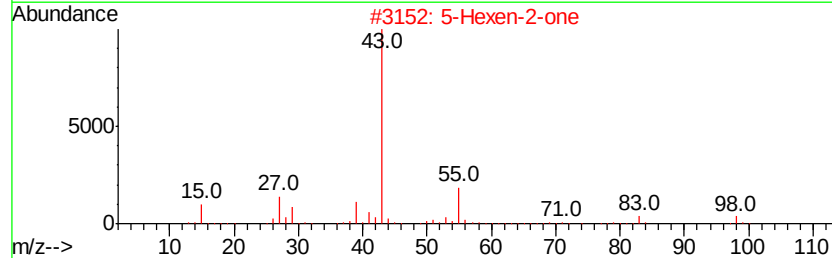
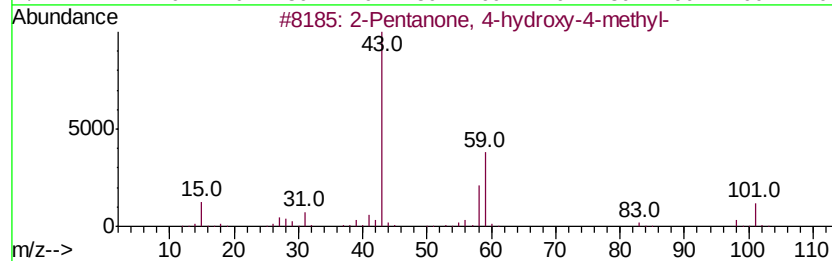
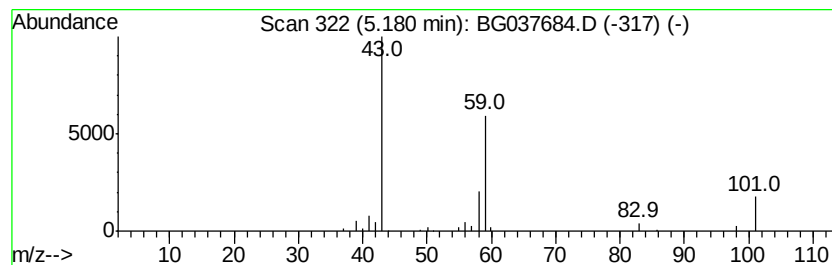
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.90 ng	46025	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9



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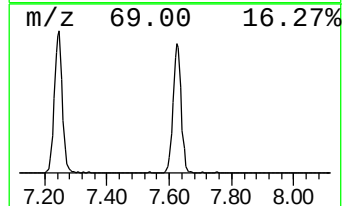
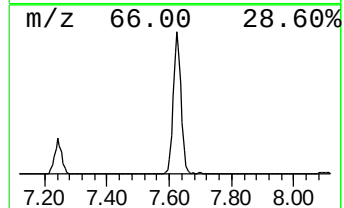
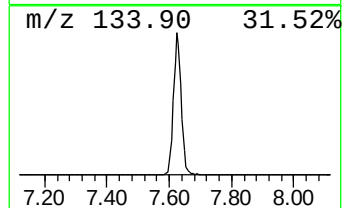
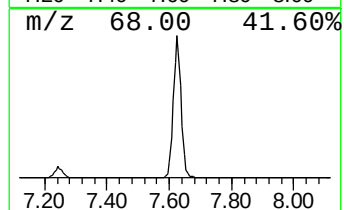
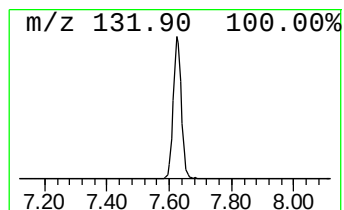
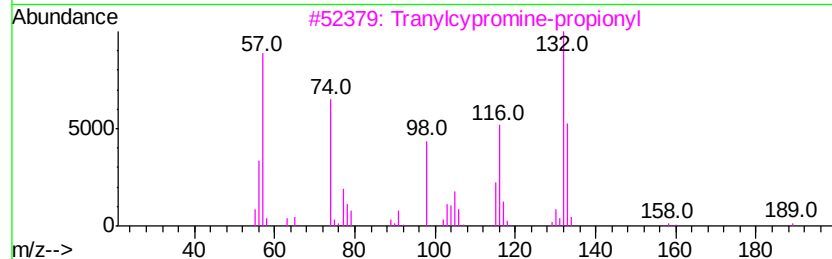
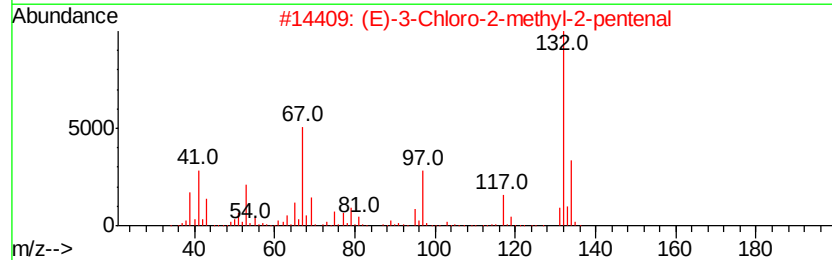
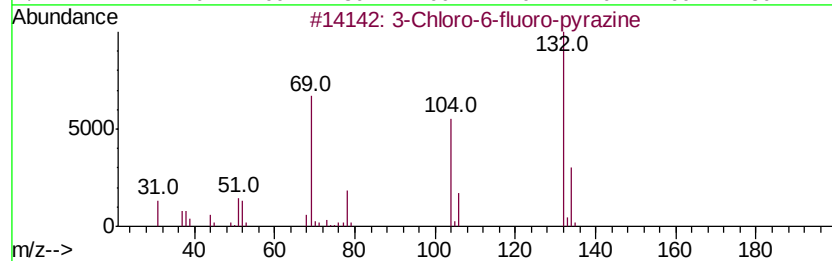
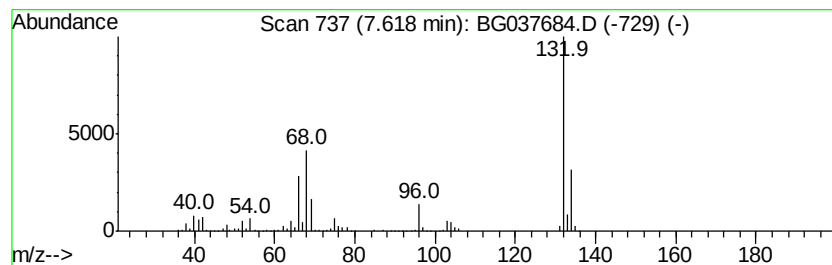
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	72.76 ng	1152940	1,4-Dichlorobenzene-d4	8.10

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	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	2.9	ng	46025	1	8.10	316919	20.0
unknown7.62	7.62	72.8	ng	1152940	1	8.10	316919	20.0