

Data Path : V:\HPCHEM1\BNA_F\DATA\BF070916\
Data File : BF088871.D
Acq On : 9 Jul 2016 15:11
Operator : UM/SJ
Sample : H3813-03 25X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P001-DL002-01

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_F\METHODS\8270-BF063016.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	1.804	18	19	34	rVB	398703	495969	8.85%	5.311%
2	3.221	108	143	144	rBV	339375	5604408	100.00%	60.013%
3	6.719	446	449	451	rBV	461839	417031	7.44%	4.466%
4	7.999	559	561	564	rBV	490545	536642	9.58%	5.746%
5	9.748	712	714	717	rBV	598190	624815	11.15%	6.691%
6	11.222	841	843	846	rBV2	549137	663850	11.85%	7.109%
7	13.851	1071	1073	1075	rBV	649019	598368	10.68%	6.407%
8	15.222	1190	1193	1196	rBV	352190	397508	7.09%	4.257%

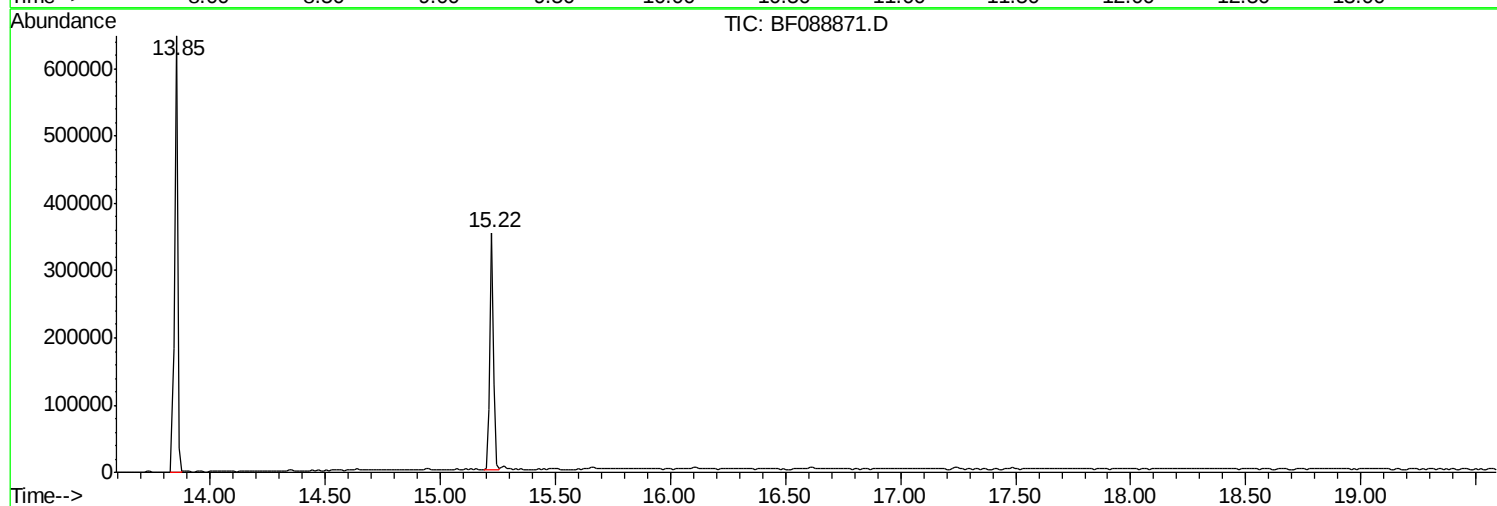
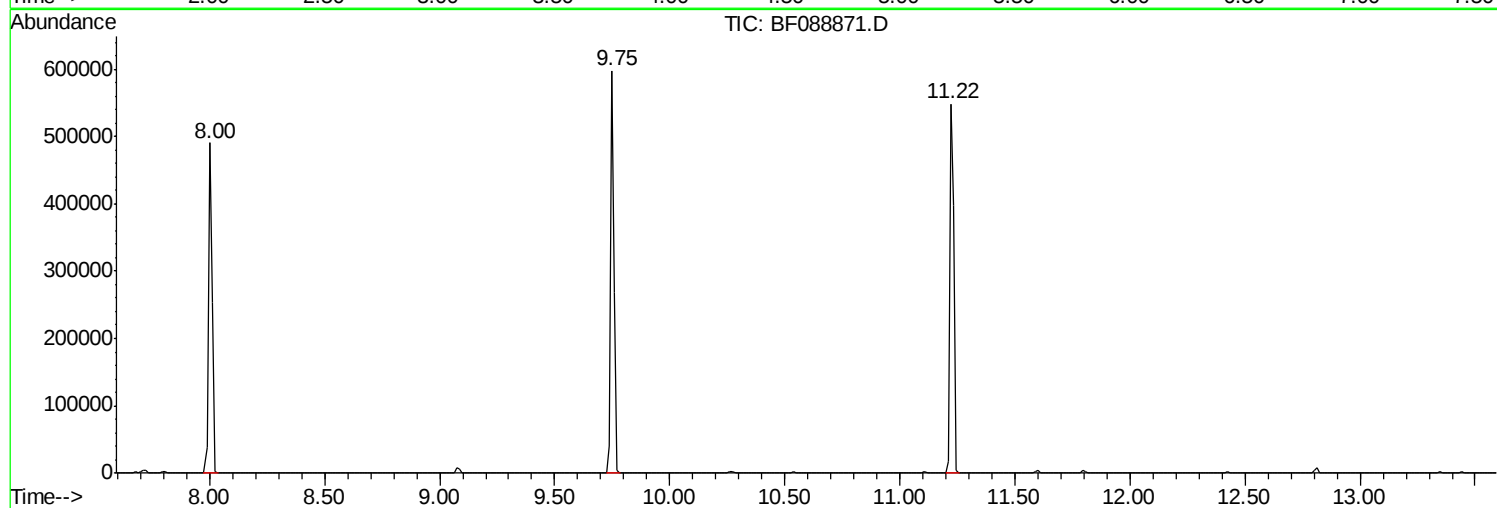
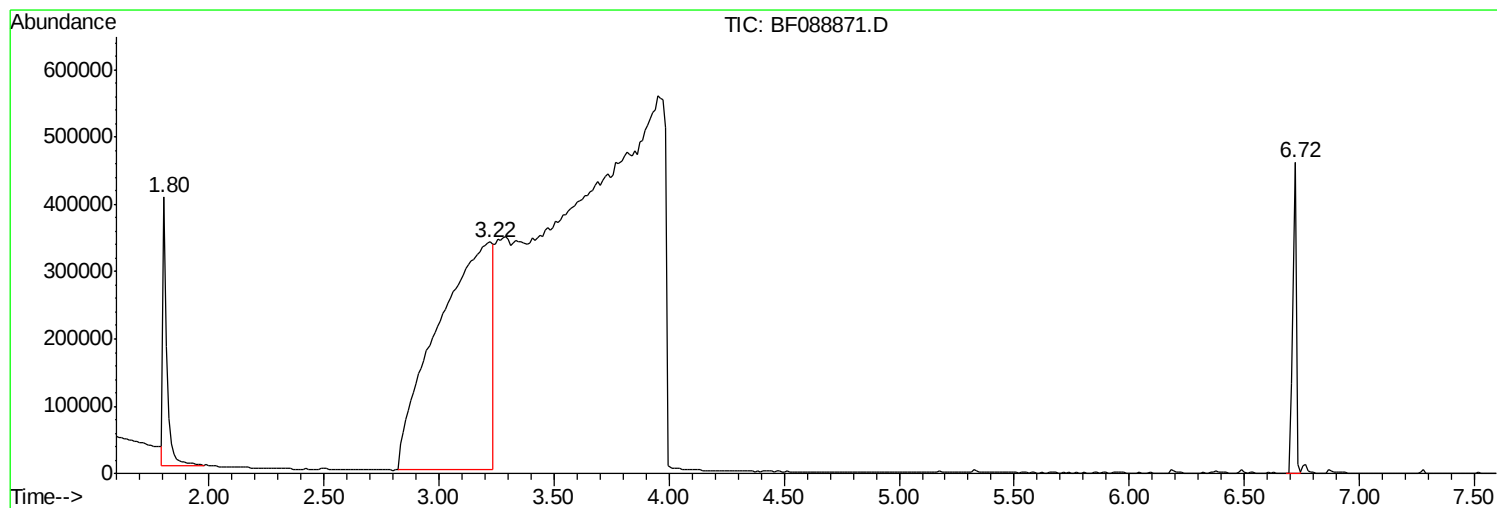
Sum of corrected areas: 9338591

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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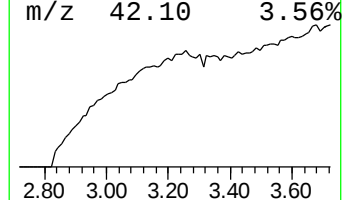
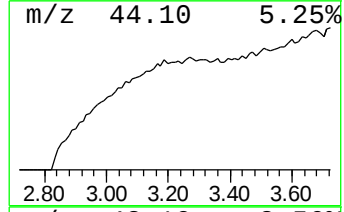
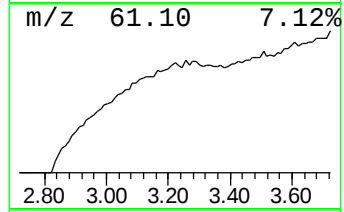
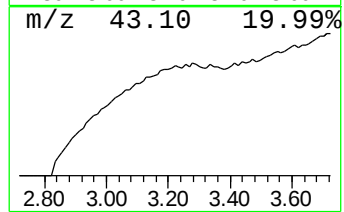
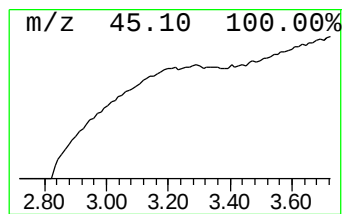
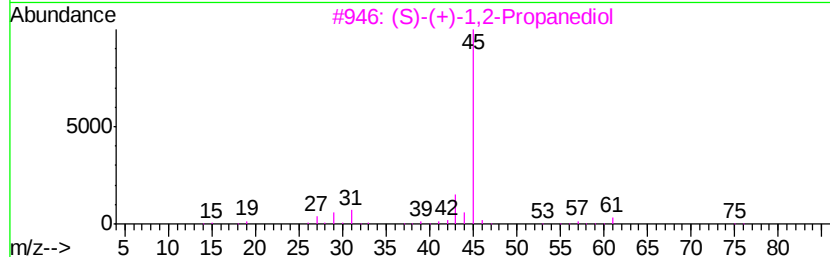
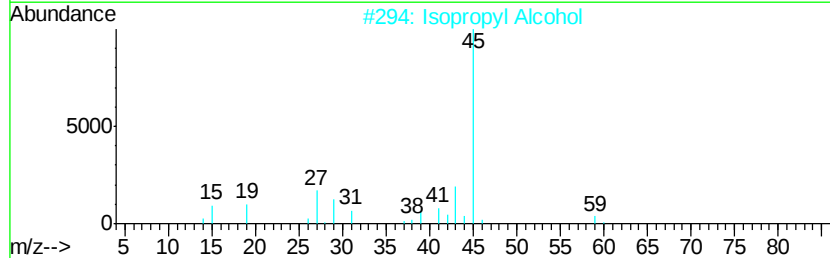
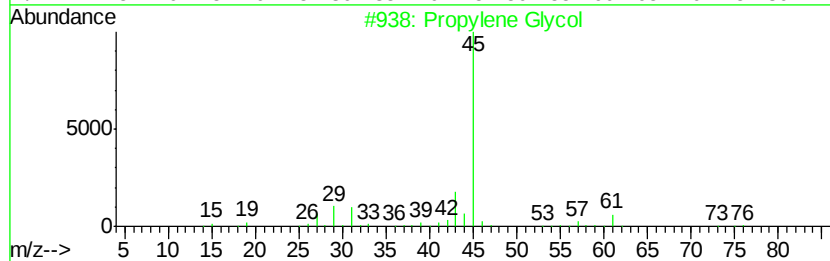
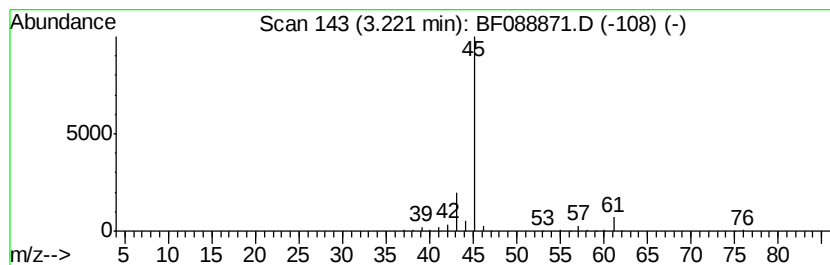
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.22	268.78 ng	5604410	1,4-Dichlorobenzene-d4	6.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propylene Glycol	76	C3H8O2	000057-55-6	90
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
4		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	9
5		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Propylene Glycol	3.22	268.8	ng	5604410	1	6.72	417031	20.0