

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072720\
 Data File : BF121026.D
 Acq On : 27 Jul 2020 11:51
 Operator : JU/CG
 Sample : PB130443BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB130443BL

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_F\METHODS\8270-BF071620.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	2.116	3	5	14	rVB	96350	109881	2.38%	0.379%
2	5.439	564	570	573	rBV	3321978	3707635	80.38%	12.788%
3	6.445	736	741	744	rBV	3099771	3595671	77.96%	12.402%
4	6.810	800	803	807	rBV	994378	900947	19.53%	3.107%
5	7.375	894	899	902	rBV	2349800	2509486	54.41%	8.656%
6	8.092	1017	1021	1025	rBV	1194888	1128002	24.46%	3.891%
7	9.175	1199	1205	1208	rBV	3874175	4344373	94.19%	14.984%
8	9.845	1315	1319	1342	rBV	1240943	1305217	28.30%	4.502%
9	10.639	1449	1454	1486	rBV	2368366	2976279	64.53%	10.266%
10	11.333	1567	1572	1580	rBV	1450616	1349316	29.25%	4.654%
11	12.921	1837	1842	1845	rBV2	4008032	4612434	100.00%	15.909%
12	13.968	2016	2020	2035	rVB	1197419	1311735	28.44%	4.524%
13	15.415	2261	2266	2290	rBV	688048	1141850	24.76%	3.938%

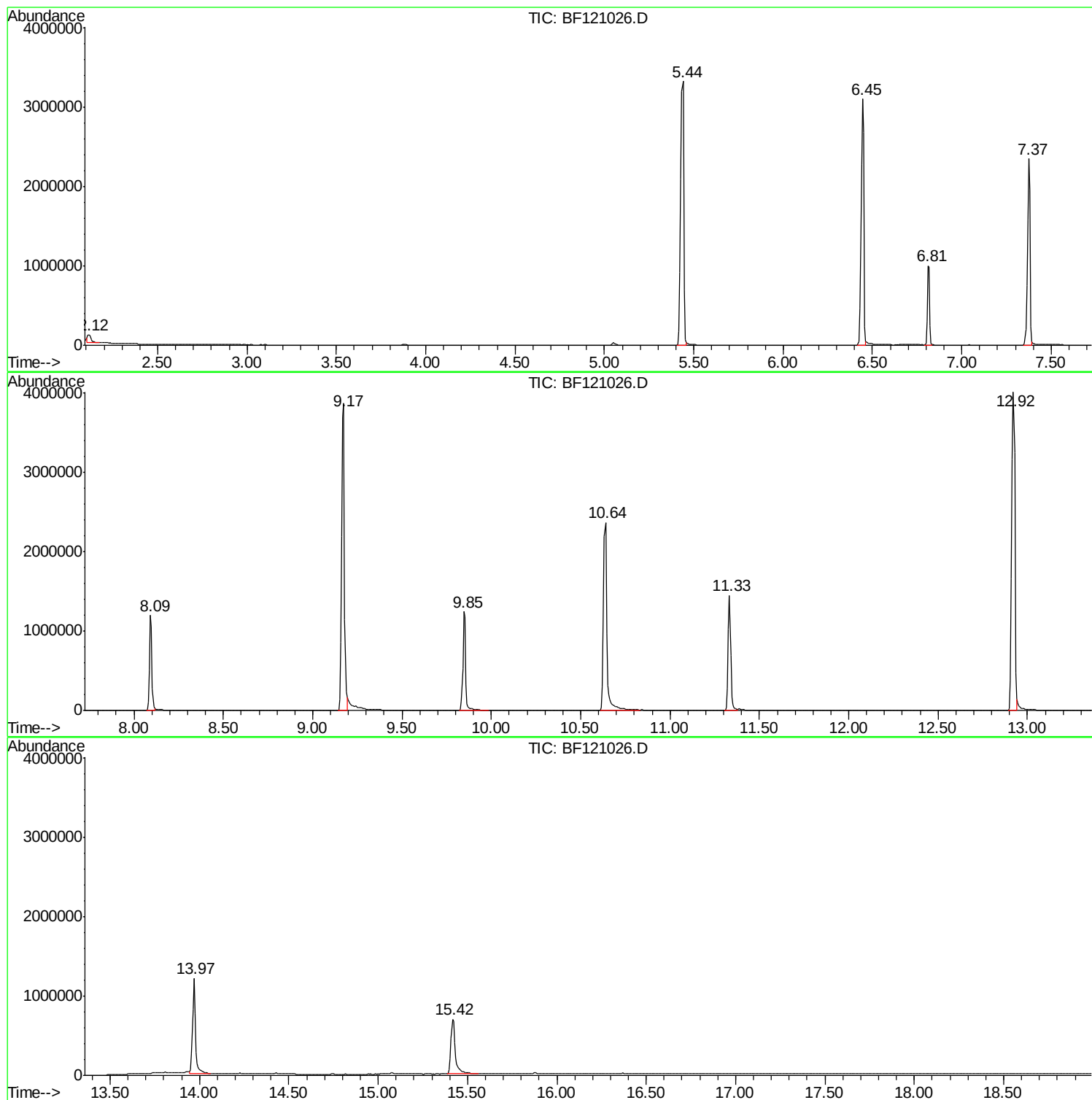
Sum of corrected areas: 28992826

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.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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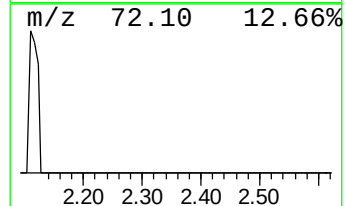
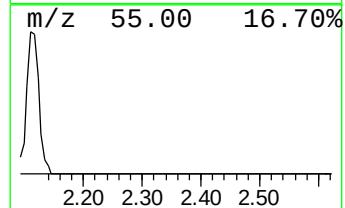
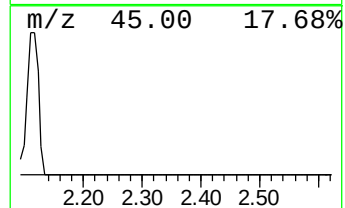
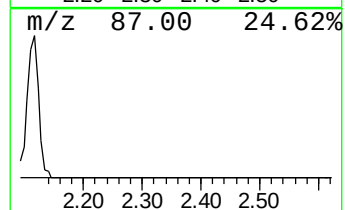
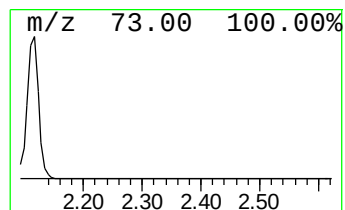
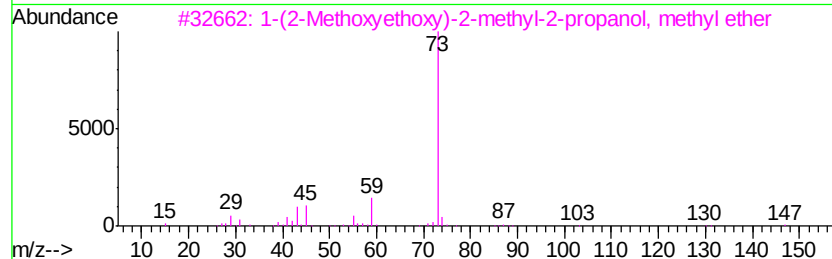
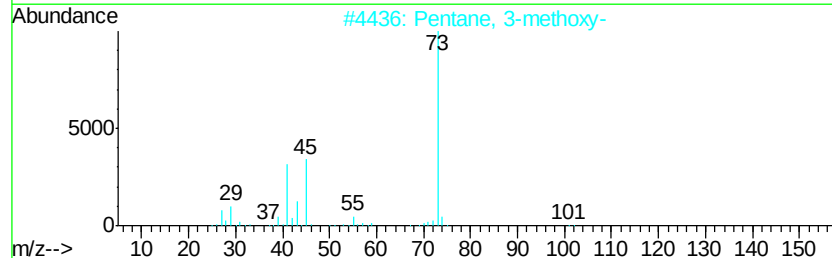
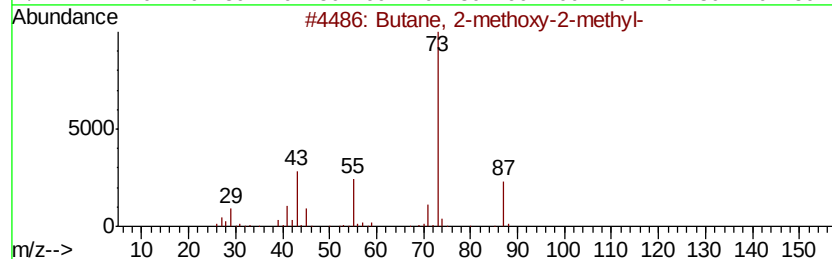
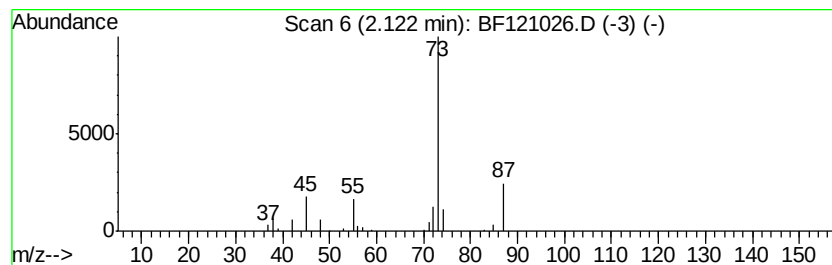
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	2.44 ng	109881	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
3		1-(2-Methoxyethoxy)-2-methyl-2-propanol, methyl ether	162	C8H18O3	1000364-24-4	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		1,3-Dioxolane, 2-pentadecyl-	284	C18H36O2	004360-57-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.12	2.4	ng	109881	1	6.82	900947	20.0