

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080119\
 Data File : BF115899.D
 Acq On : 1 Aug 2019 11:22
 Operator : HP/JU
 Sample : PB121853BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB121853BL

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-BF073019.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	2.199	10	19	26	rVB	83699	121800	2.36%	0.327%
2	5.487	573	578	581	rBV	3467679	3544399	68.63%	9.516%
3	6.492	743	749	752	rBV	3415529	3725031	72.13%	10.001%
4	6.628	767	772	775	rBV	3808999	3851009	74.57%	10.339%
5	6.851	807	810	820	rVB	1112935	955848	18.51%	2.566%
6	7.010	833	837	841	rBV	3311639	3121217	60.43%	8.380%
7	7.416	900	906	909	rBV	2404169	2495995	48.33%	6.701%
8	8.134	1024	1028	1048	rBV	1344497	1224263	23.70%	3.287%
9	9.216	1206	1212	1215	rBV	4610270	4641420	89.87%	12.461%
10	9.892	1322	1327	1341	rBV	1583713	1483378	28.72%	3.983%
11	10.686	1457	1462	1490	rBV	2800054	2972182	57.55%	7.980%
12	11.380	1576	1580	1588	rBV	1668391	1525061	29.53%	4.094%
13	12.969	1845	1850	1853	rBV	4559261	5164609	100.00%	13.866%
14	14.021	2025	2029	2042	rBV	1308864	1287692	24.93%	3.457%
15	15.498	2274	2280	2293	rBV	811364	1133470	21.95%	3.043%

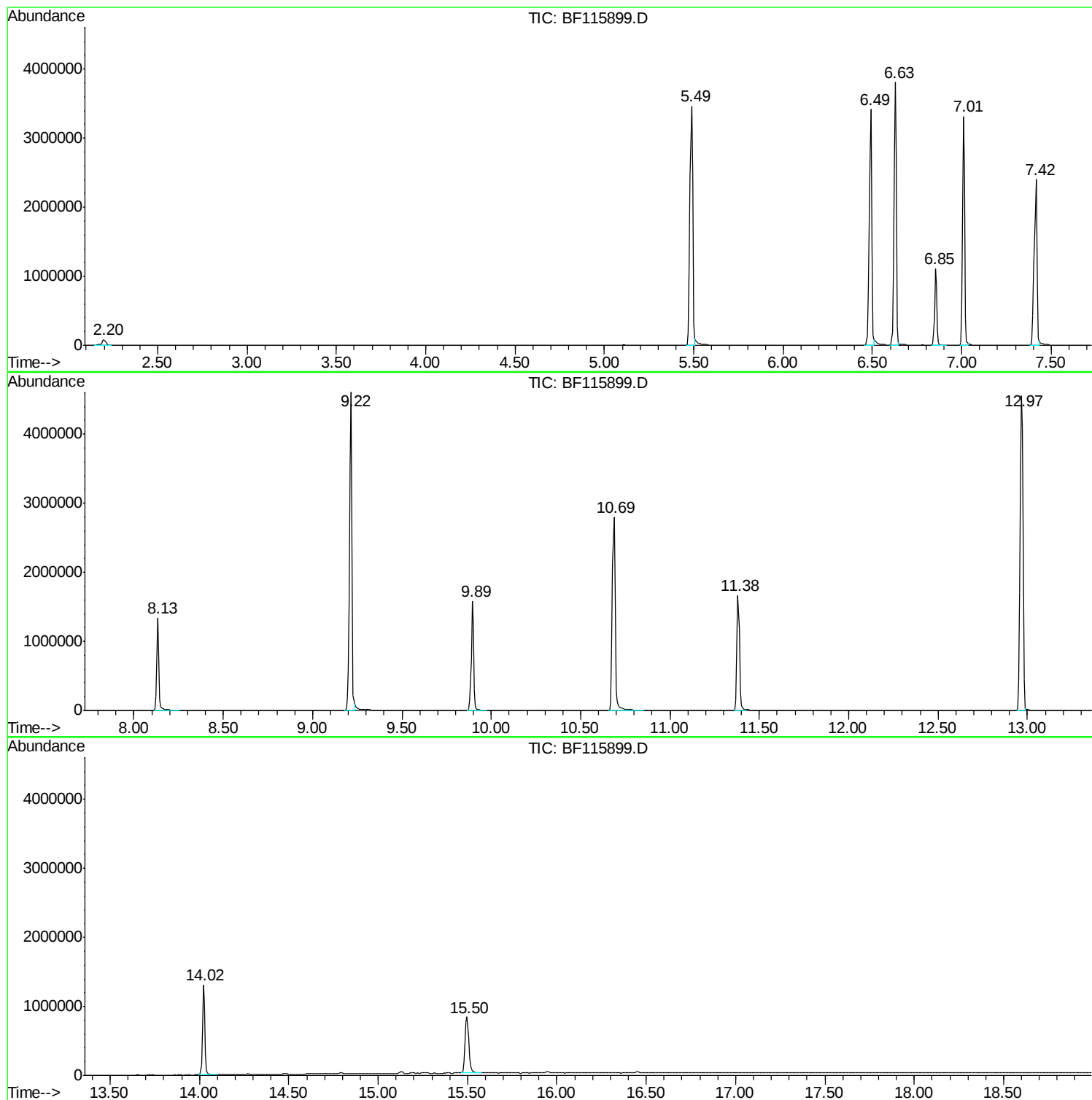
Sum of corrected areas: 37247374

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



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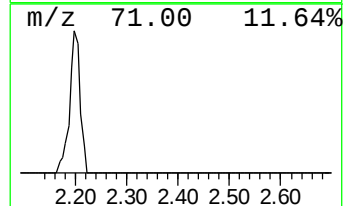
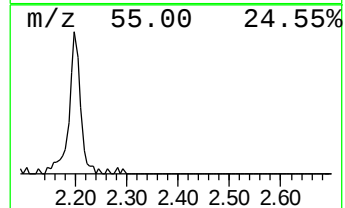
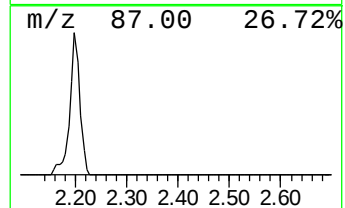
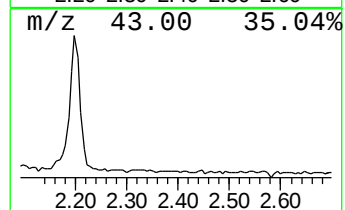
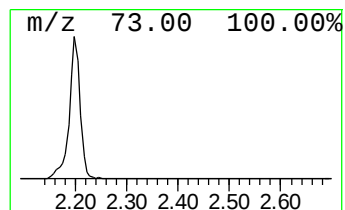
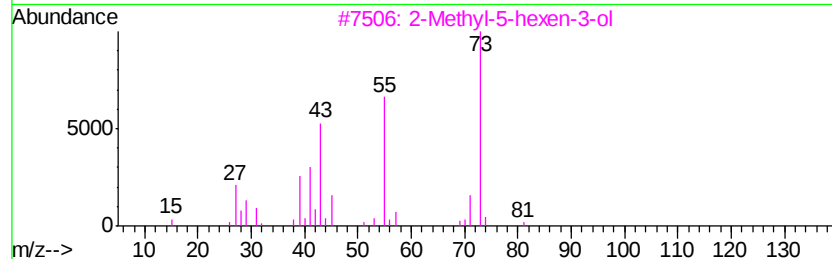
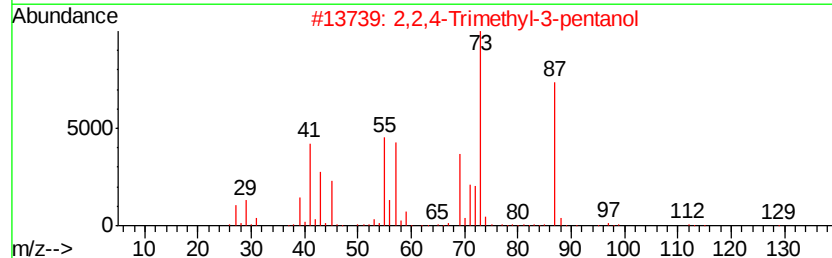
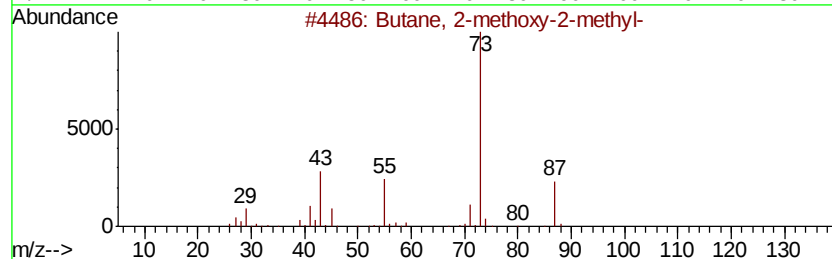
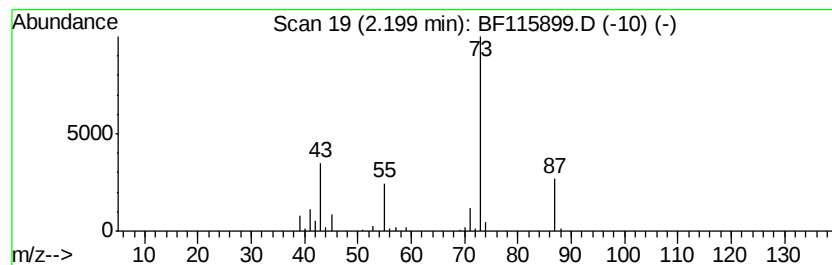
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.20	2.55 ng	121800	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17



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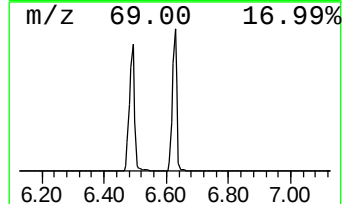
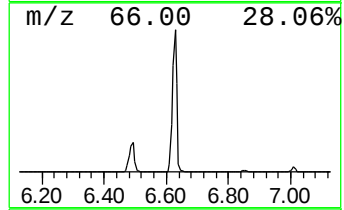
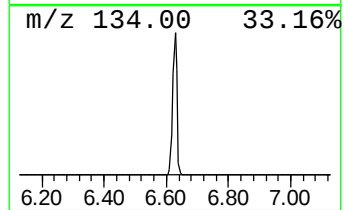
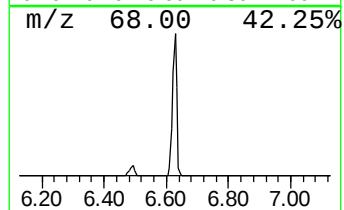
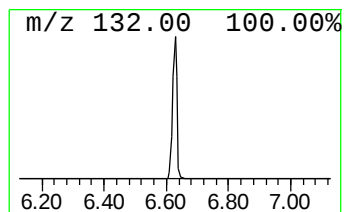
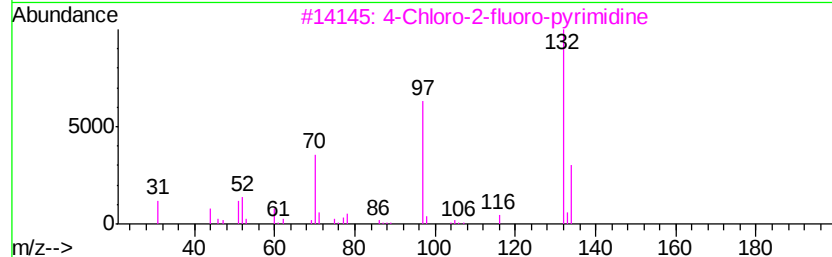
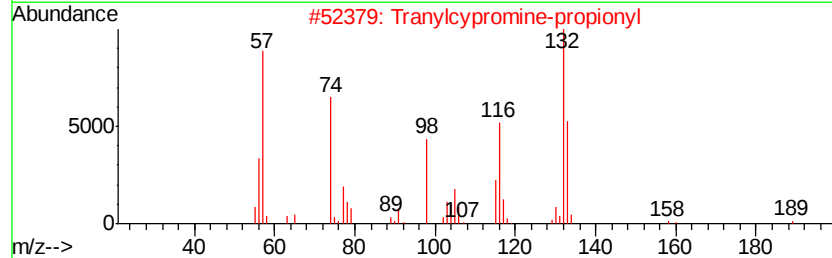
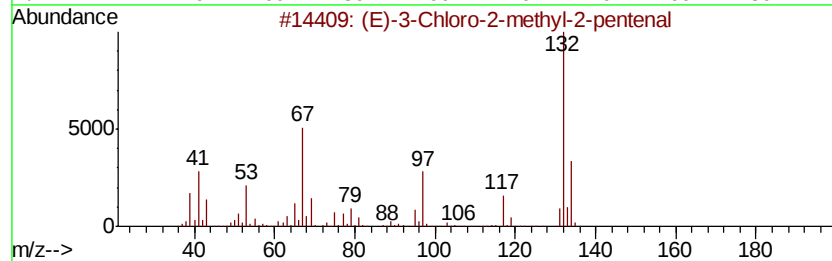
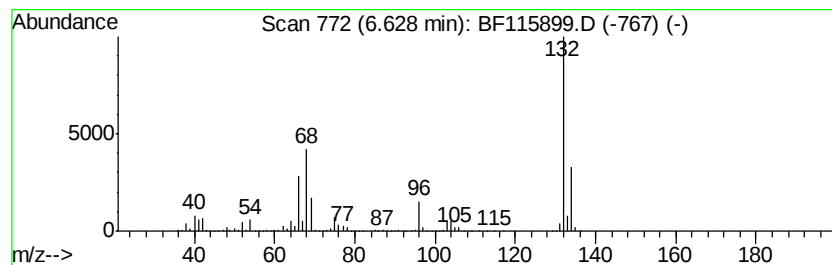
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	80.58 ng	3851010	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
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
Butane, 2-methoxy...	2.20	2.5	ng	121800	1	6.85	955848	20.0
unknown6.63	6.63	80.6	ng	3851010	1	6.85	955848	20.0