

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071819\
 Data File : BF115646.D
 Acq On : 18 Jul 2019 14:05
 Operator : HP/JU
 Sample : PB121479BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB121479BL

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-BF071219.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	2.299	31	36	44	rVB	202182	289583	4.72%	0.614%
2	5.522	578	584	587	rBV	4254782	4701358	76.70%	9.972%
3	6.516	747	753	756	rBV	3931365	4932174	80.47%	10.462%
4	6.657	772	777	780	rBV	4500232	5027324	82.02%	10.664%
5	6.881	812	815	819	rBV	1354183	1186752	19.36%	2.517%
6	7.040	837	842	846	rBV	3765535	4025455	65.68%	8.539%
7	7.445	904	911	914	rBV	2831113	3318046	54.13%	7.038%
8	8.163	1028	1033	1037	rBV	1674500	1550842	25.30%	3.290%
9	9.239	1209	1216	1219	rBV	5617917	6129342	100.00%	13.001%
10	9.916	1325	1331	1344	rBV	2140527	1880369	30.68%	3.989%
11	10.704	1459	1465	1491	rBV	3522891	3849167	62.80%	8.165%
12	11.398	1579	1583	1587	rBV	2107524	1875707	30.60%	3.979%
13	12.986	1848	1853	1856	rBV	5600002	5964914	97.32%	12.652%
14	14.033	2027	2031	2040	rBV	1444460	1322181	21.57%	2.805%
15	15.498	2275	2280	2293	rBV2	796845	1091068	17.80%	2.314%

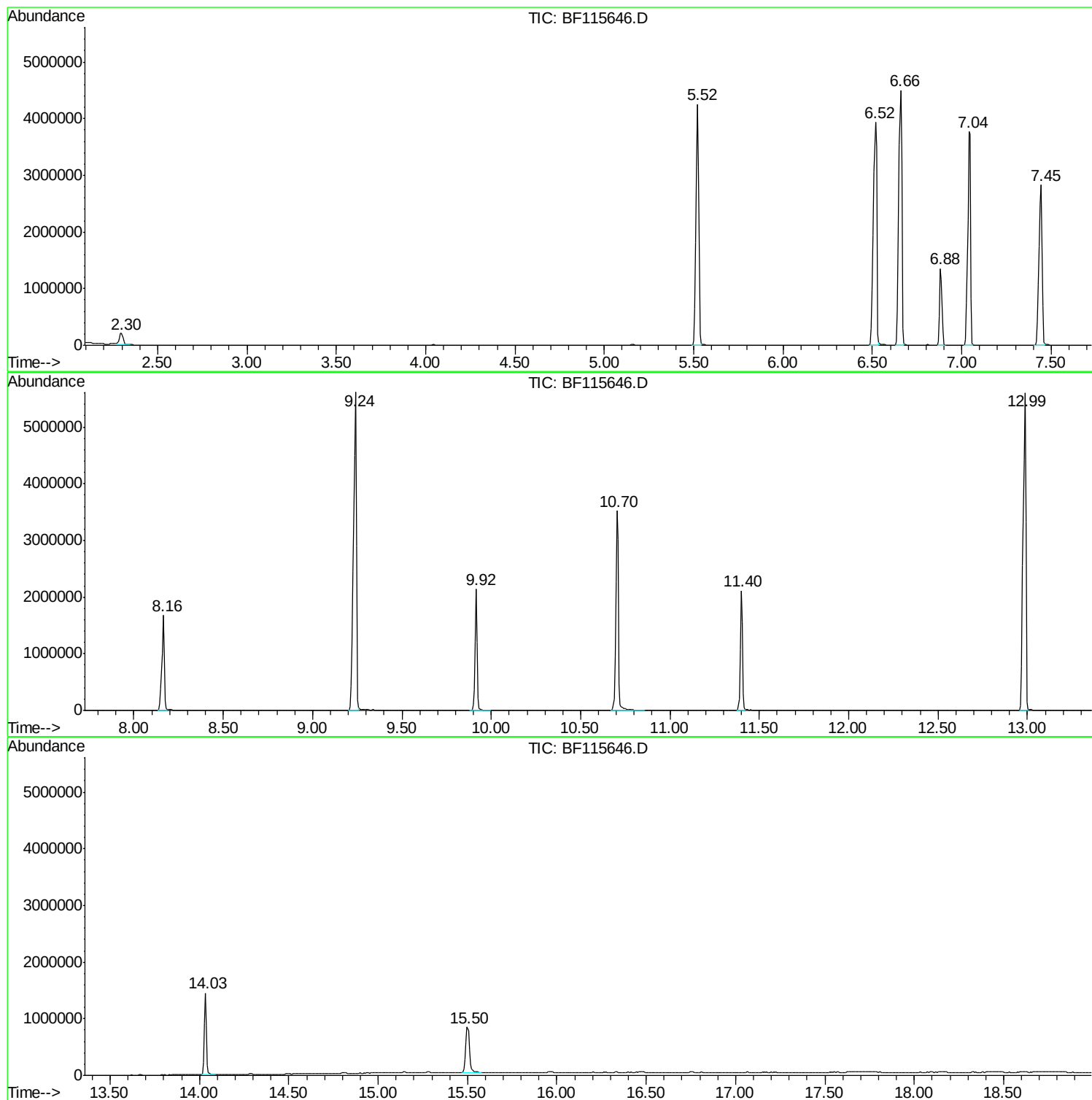
Sum of corrected areas: 47144282

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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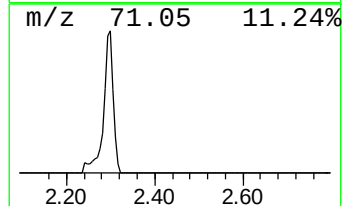
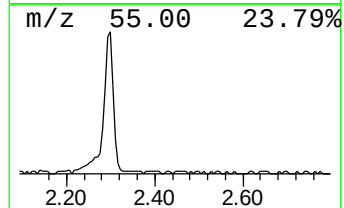
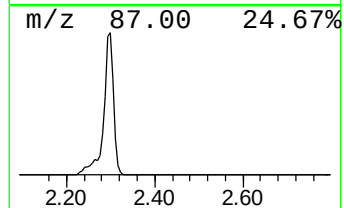
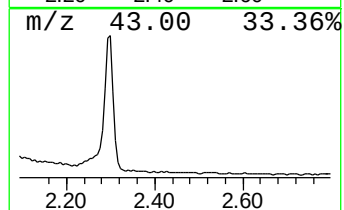
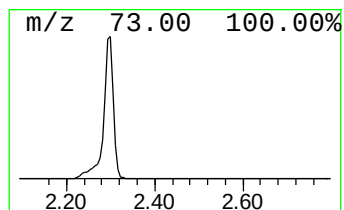
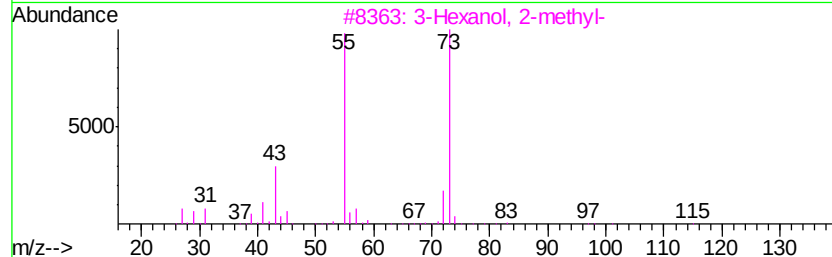
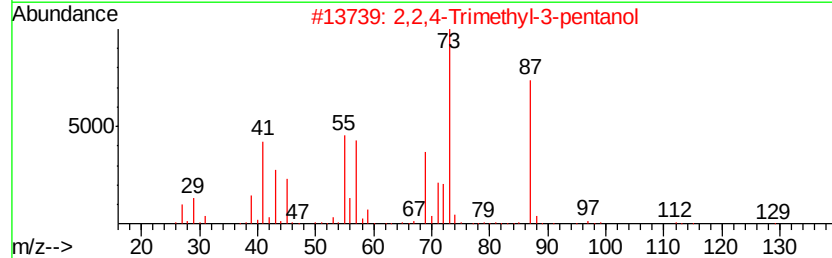
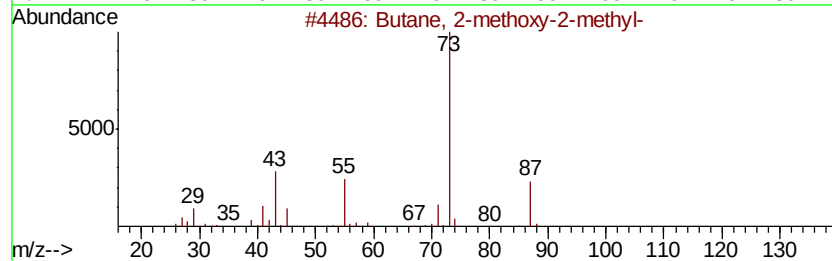
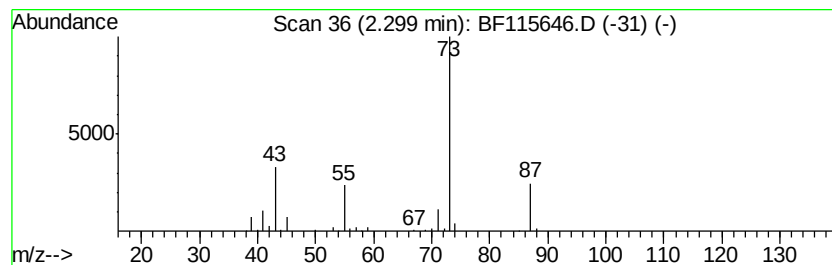
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.30	4.88 ng	289583	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9



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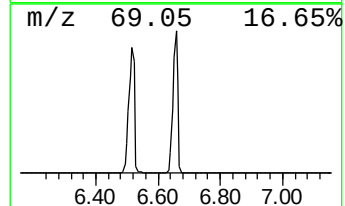
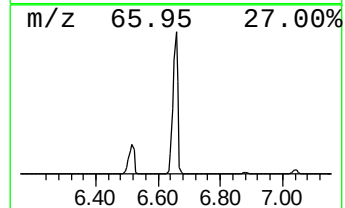
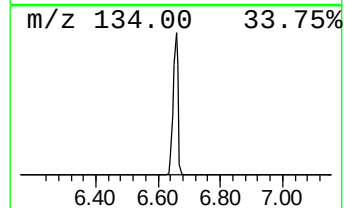
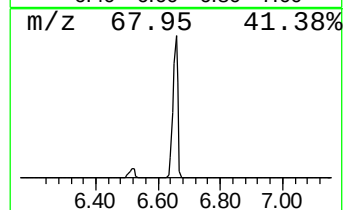
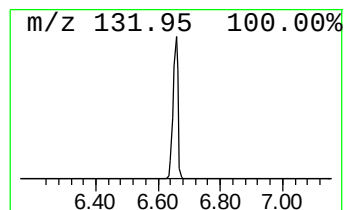
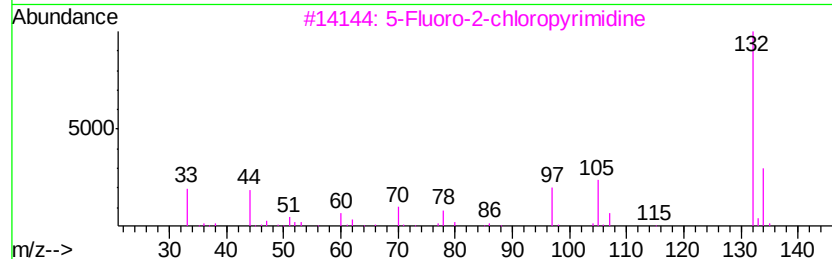
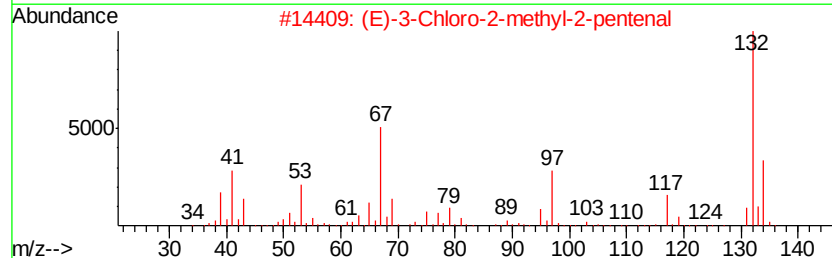
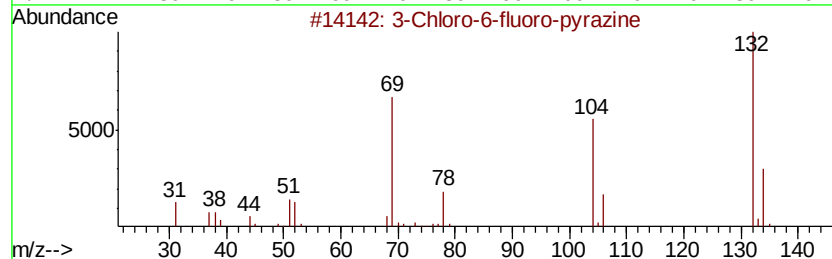
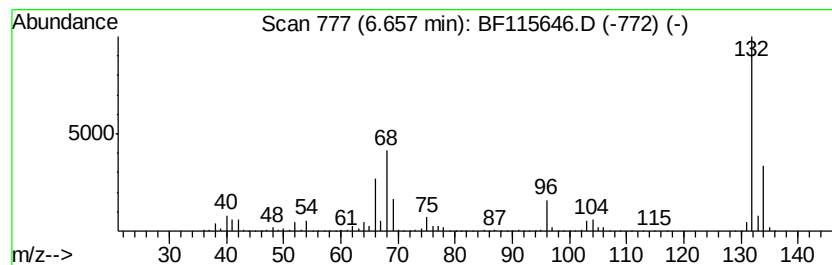
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	84.72 ng	5027320	1,4-Dichlorobenzene-d4	6.88

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	12
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.30	4.9	ng	289583	1	6.88	1186750	20.0
unknown6.66	6.66	84.7	ng	5027320	1	6.88	1186750	20.0