

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
 Data File : BF113963.D
 Acq On : 24 Apr 2019 16:03
 Operator : JU/SJ
 Sample : K2487-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB

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-BF042219.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	2.193	14	18	43	rVB	184745	380591	8.32%	1.312%
2	5.516	579	583	586	rBV	1742171	1567090	34.24%	5.403%
3	6.516	749	753	766	rBV	1216952	1029824	22.50%	3.550%
4	6.663	774	778	782	rBV	3045672	3045514	66.55%	10.499%
5	6.892	814	817	821	rBV	904347	787981	17.22%	2.717%
6	7.051	840	844	848	rBV	3052582	2914993	63.70%	10.050%
7	7.457	907	913	916	rBV	2208929	2279144	49.80%	7.857%
8	8.175	1031	1035	1048	rBV	1099123	972880	21.26%	3.354%
9	9.251	1213	1218	1221	rBV	4800742	4408334	96.33%	15.198%
10	9.927	1330	1333	1348	rVB	1178902	1151763	25.17%	3.971%
11	10.716	1463	1467	1483	rBV	2639594	2615021	57.14%	9.015%
12	11.416	1582	1586	1594	rBV	1288864	1186601	25.93%	4.091%
13	12.998	1850	1855	1858	rBV	4594619	4576359	100.00%	15.777%
14	14.057	2031	2035	2042	rVB	1063258	1005014	21.96%	3.465%
15	15.562	2285	2291	2296	rBV	799280	1085176	23.71%	3.741%

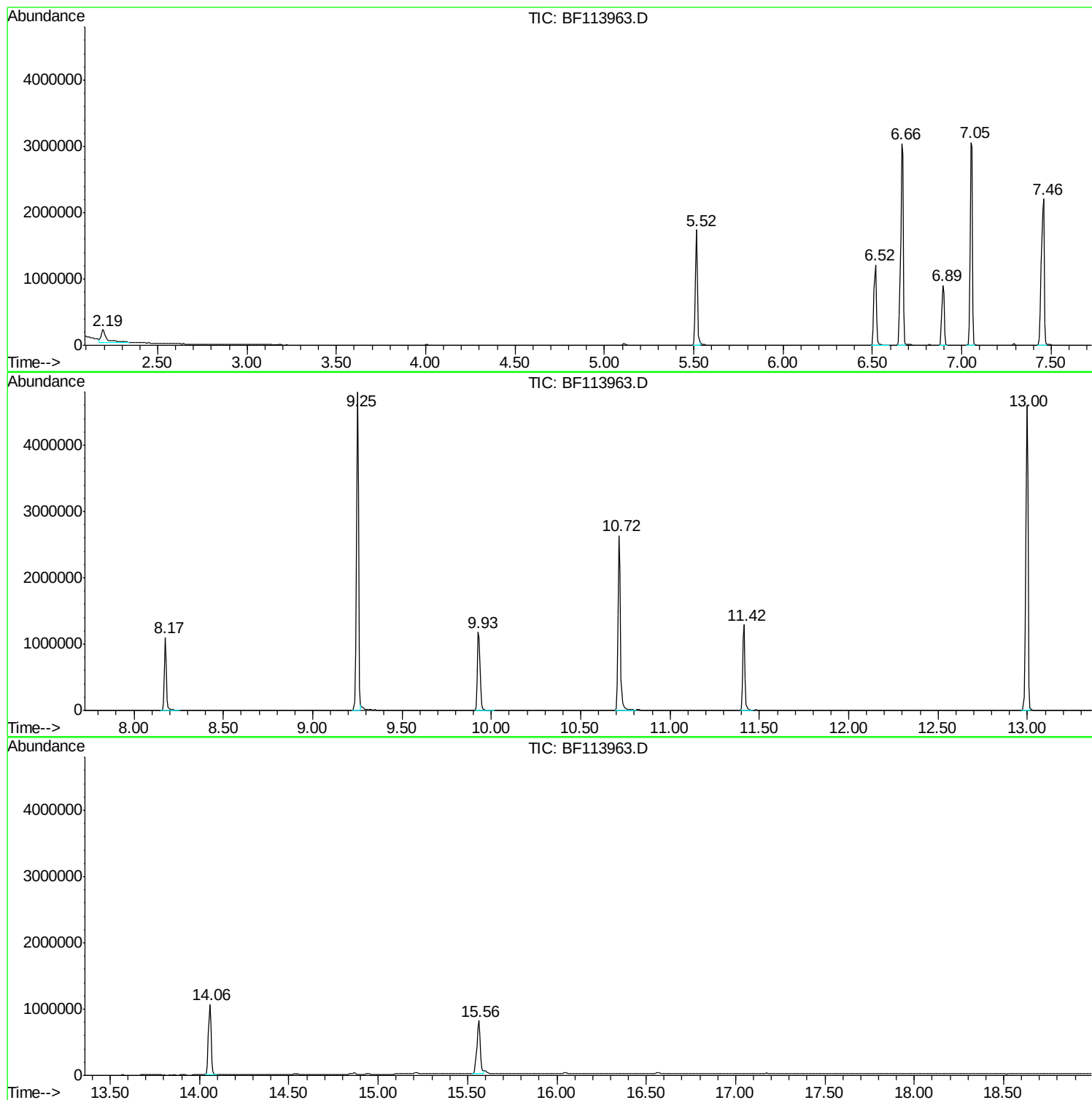
Sum of corrected areas: 29006285

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.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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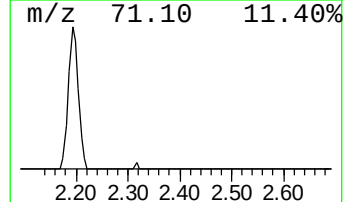
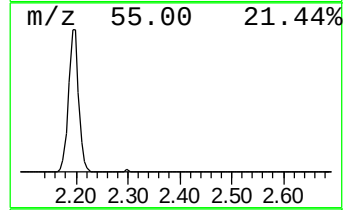
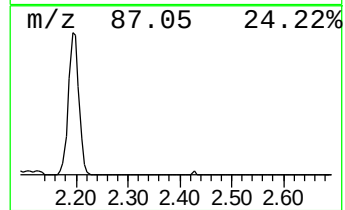
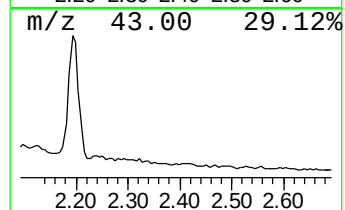
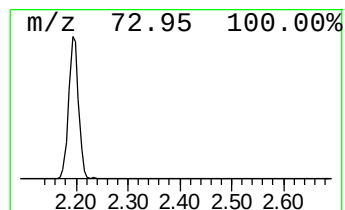
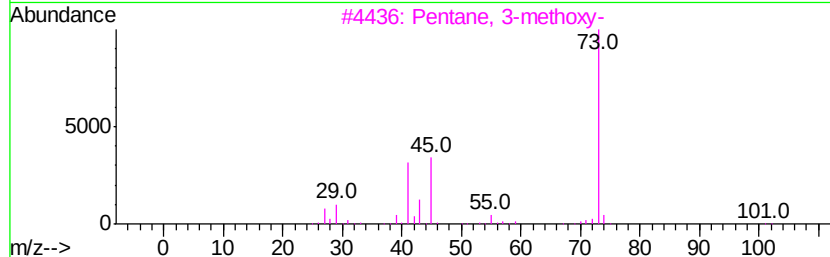
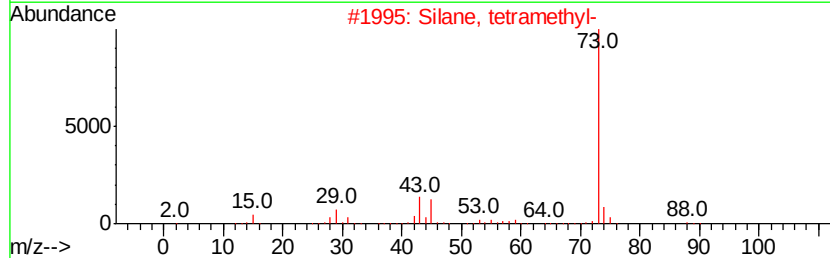
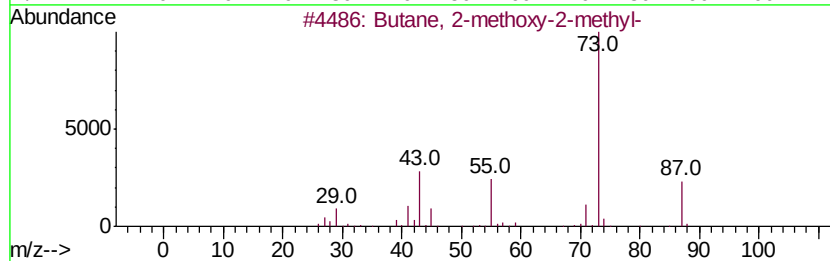
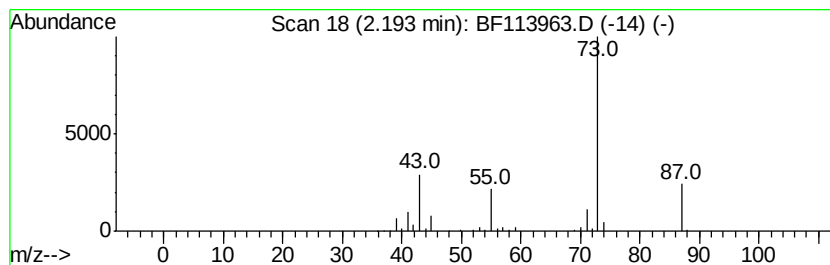
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	9.66 ng	380591	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5			Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	12



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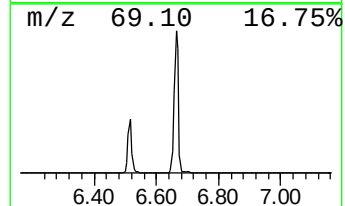
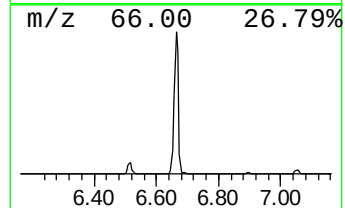
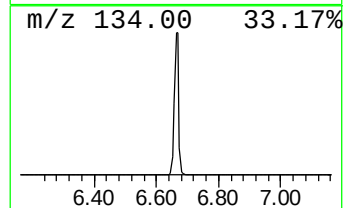
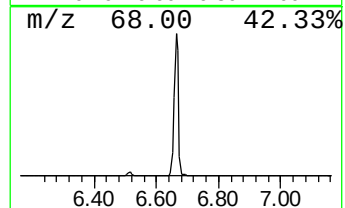
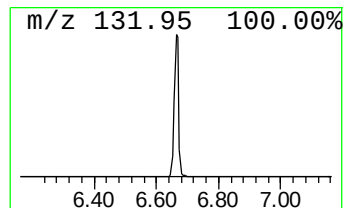
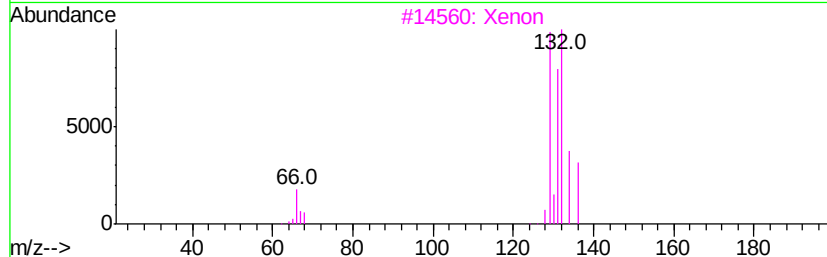
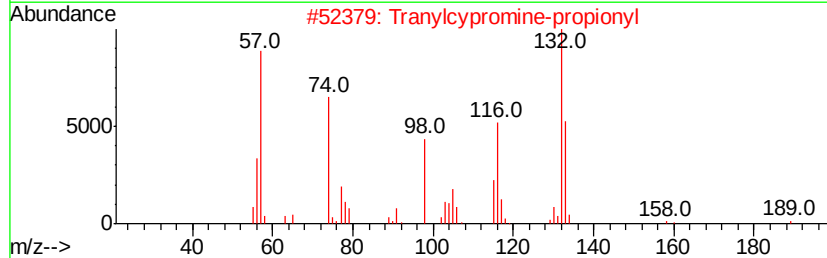
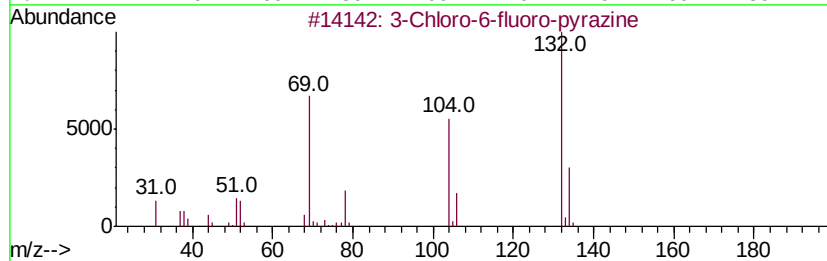
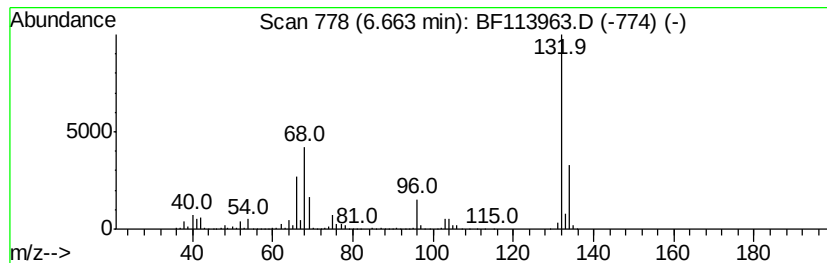
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	77.30 ng	3045510	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Xenon	132	Xe	007440-63-3	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.19	9.7	ng	380591	1	6.89	787981	20.0
unknown6.66	6.66	77.3	ng	3045510	1	6.89	787981	20.0