

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081519\
 Data File : BF116274.D
 Acq On : 15 Aug 2019 12:56
 Operator : HP/JU
 Sample : PB122172BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122172BL

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	2.146	6	10	19	rVB	94665	140865	2.84%	0.385%
2	5.451	568	572	602	rBV	3283322	3583439	72.33%	9.783%
3	6.463	739	744	762	rBV	3131706	3736535	75.42%	10.201%
4	6.593	762	766	769	rBV	3773078	3639216	73.46%	9.935%
5	6.816	801	804	817	rVB	967029	885888	17.88%	2.419%
6	6.975	827	831	848	rBV	3186728	3079115	62.15%	8.406%
7	7.381	896	900	924	rBV	2294103	2558695	51.65%	6.985%
8	8.098	1018	1022	1041	rBV	1123734	1146662	23.15%	3.130%
9	9.169	1200	1204	1218	rBV	4090815	4571285	92.27%	12.480%
10	9.851	1316	1320	1336	rBV	1365557	1384165	27.94%	3.779%
11	10.639	1450	1454	1483	rBV	2577111	2922976	59.00%	7.980%
12	11.333	1569	1572	1594	rBV	1285764	1441097	29.09%	3.934%
13	12.922	1837	1842	1845	rBV	4604060	4954242	100.00%	13.526%
14	13.974	2018	2021	2037	rBV	1023590	1236949	24.97%	3.377%
15	15.051	2199	2204	2211	rBV	59418	73431	1.48%	0.200%
16	15.433	2262	2269	2290	rVB	706880	1134607	22.90%	3.098%
17	15.839	2333	2338	2348	rVB	49225	80508	1.63%	0.220%
18	16.327	2415	2421	2432	rVB	28826	59136	1.19%	0.161%

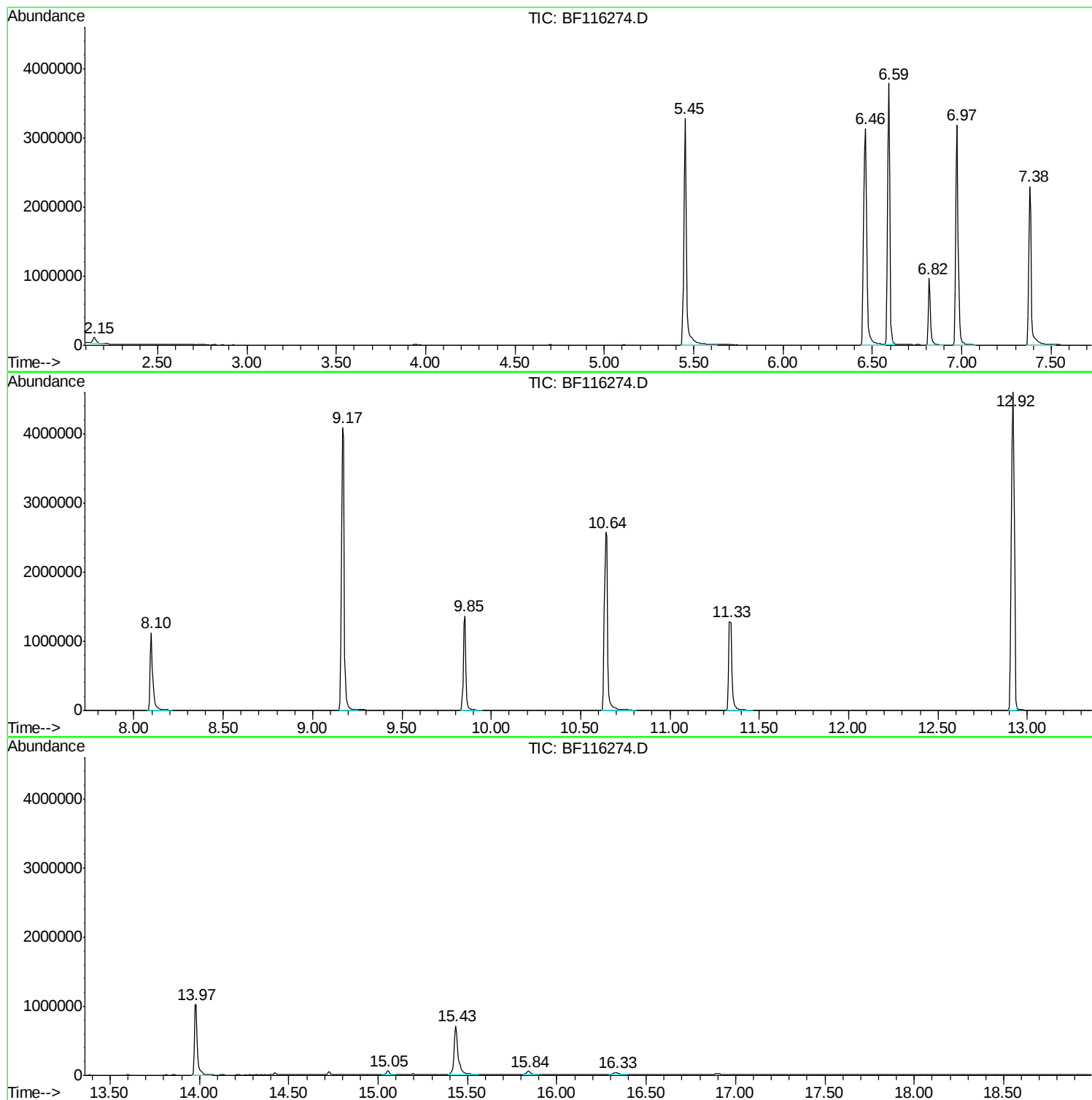
Sum of corrected areas: 36628811

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.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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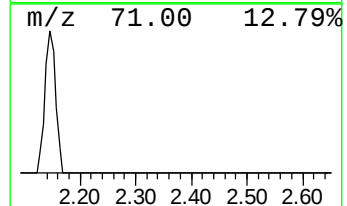
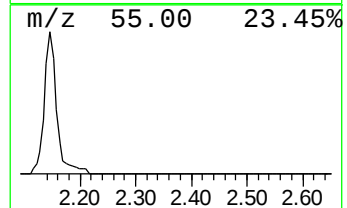
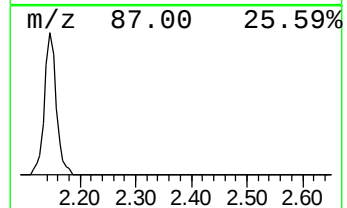
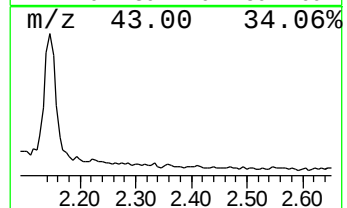
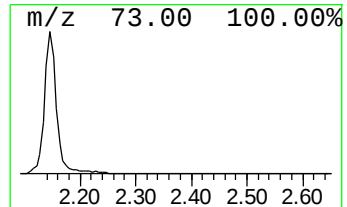
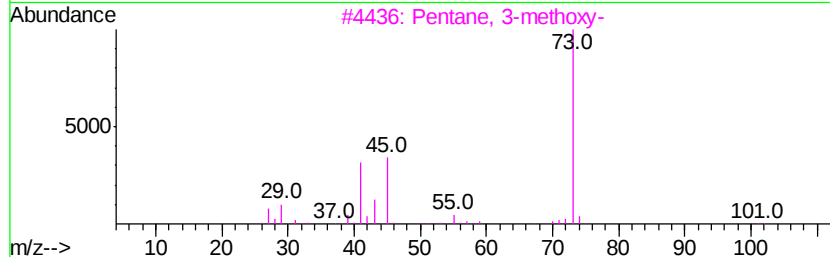
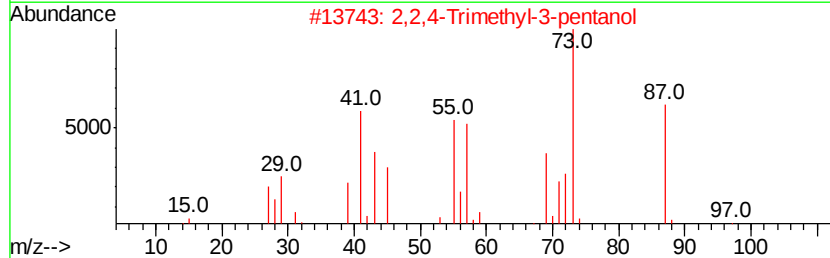
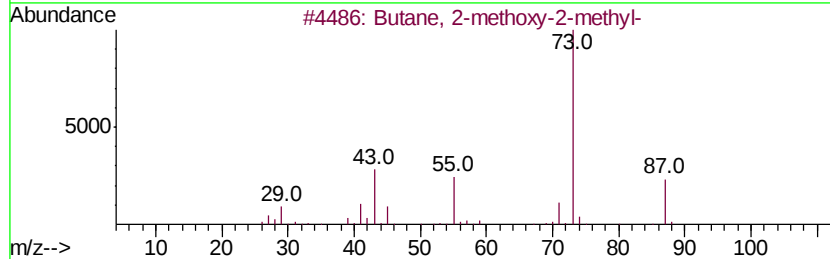
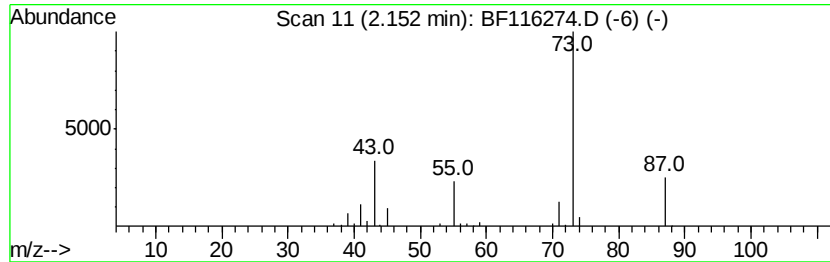
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
2.15	3.18 ng	140865	1,4-Dichlorobenzene-d4		6.82	
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	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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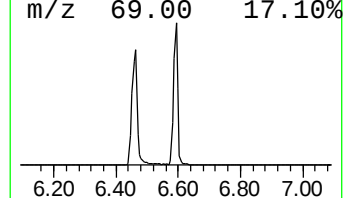
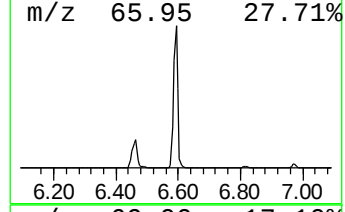
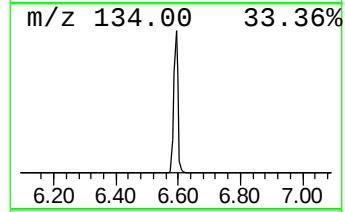
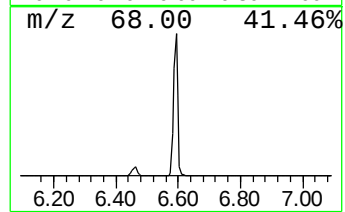
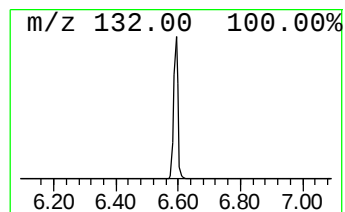
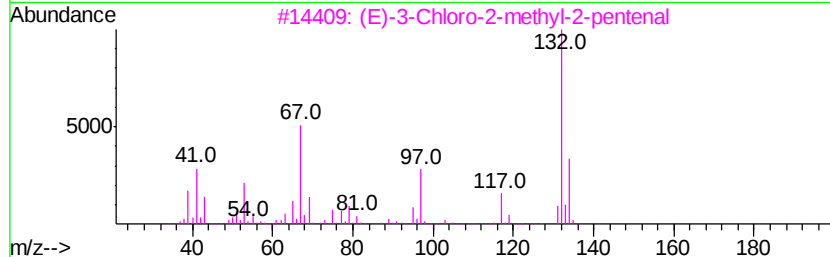
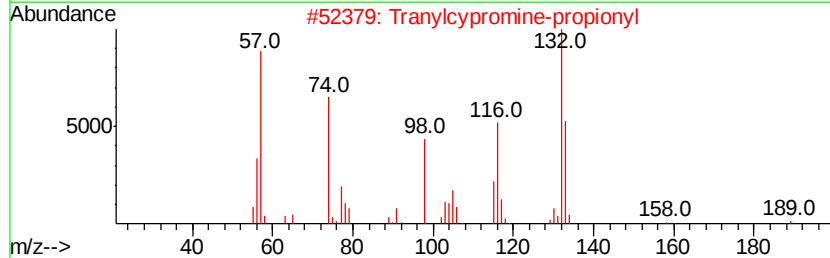
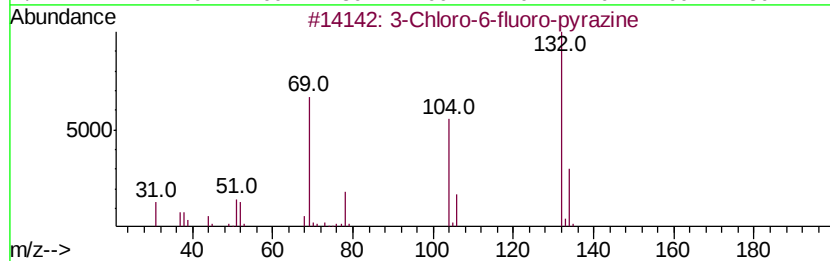
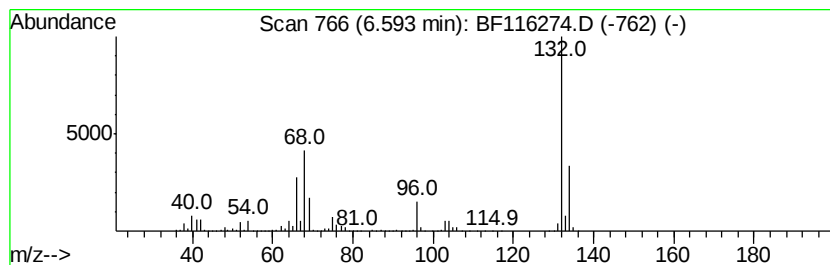
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Peak Number 2 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	82.16 ng	3639220	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
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...	2.15	3.2	ng	140865	1	6.82	885888	20.0
unknown6.59	6.59	82.2	ng	3639220	1	6.82	885888	20.0