

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102424\
 Data File : BF139993.D
 Acq On : 24 Oct 2024 11:20
 Operator : RC/JU
 Sample : PB164315BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164315BL

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-BF101824.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139993.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.240	14	25	33	rVB	88056	160889	2.95%	0.477%
2	5.134	508	517	527	rBV	200198	338596	6.20%	1.004%
3	5.516	575	582	587	rBV	3239613	4596827	84.23%	13.630%
4	6.510	744	751	756	rBV	3053922	4521906	82.85%	13.408%
5	6.886	810	815	821	rBV	716016	898847	16.47%	2.665%
6	7.451	904	911	916	rBV	2090918	3070735	56.26%	9.105%
7	8.169	1027	1033	1038	rBV	922603	1186900	21.75%	3.519%
8	9.245	1209	1216	1221	rBV	3932333	5445446	99.77%	16.146%
9	9.927	1325	1332	1344	rBV	1023782	1370406	25.11%	4.063%
10	10.716	1459	1466	1488	rBV	2571227	3510641	64.32%	10.409%
11	11.410	1579	1584	1590	rBV	1106437	1417446	25.97%	4.203%
12	13.004	1848	1855	1860	rBV	3885327	5457759	100.00%	16.183%
13	14.051	2027	2033	2045	rBV	756222	988364	18.11%	2.931%
14	15.527	2278	2284	2300	rBV	441477	760882	13.94%	2.256%

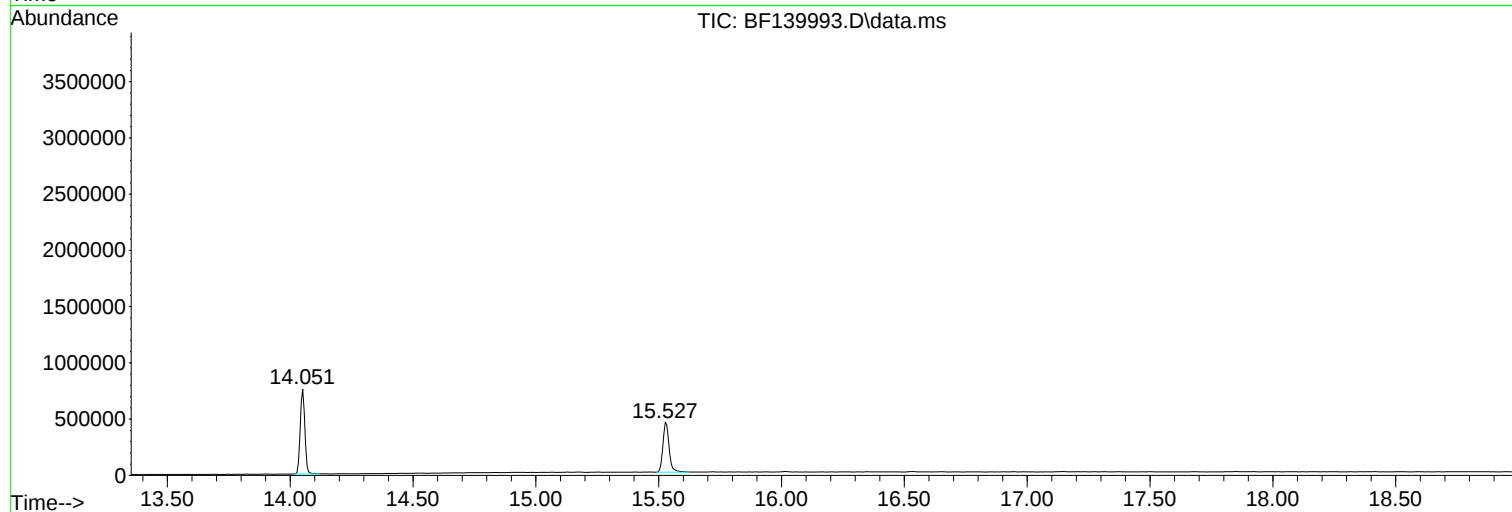
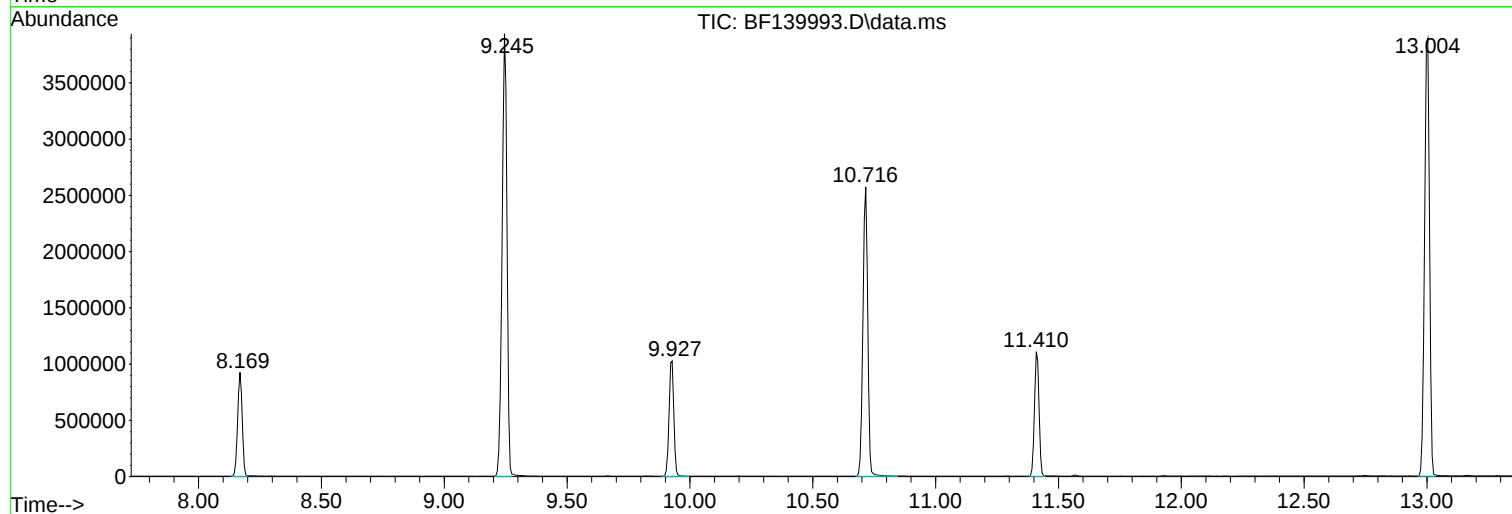
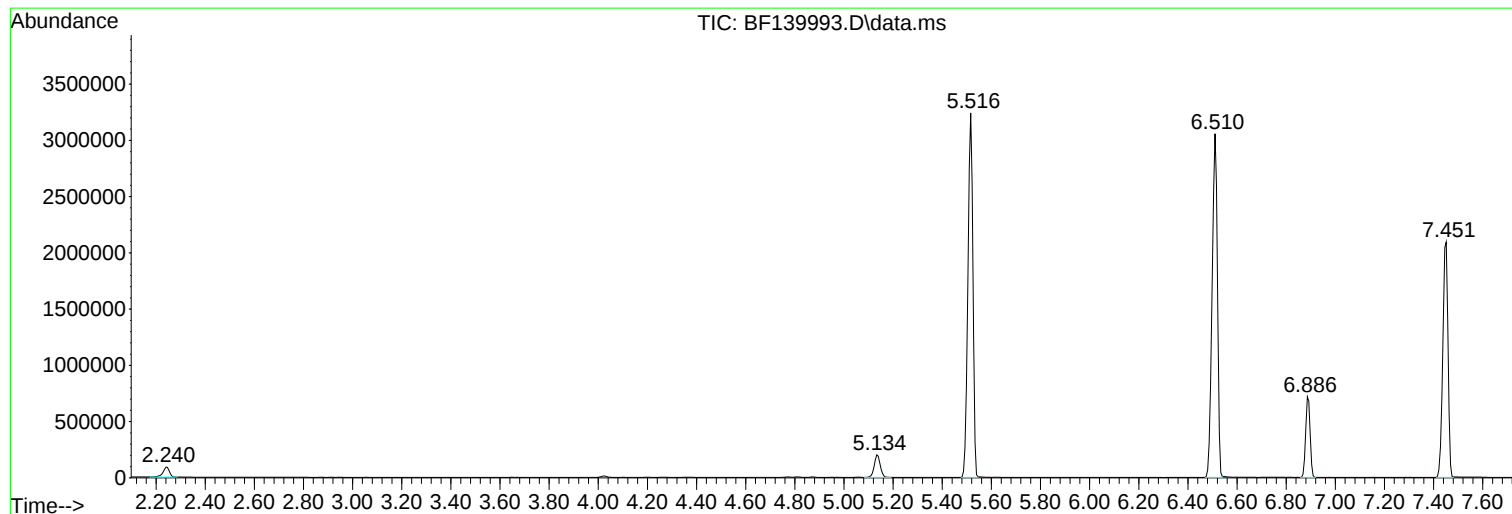
Sum of corrected areas: 33725644

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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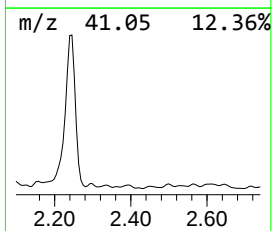
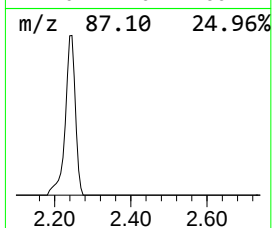
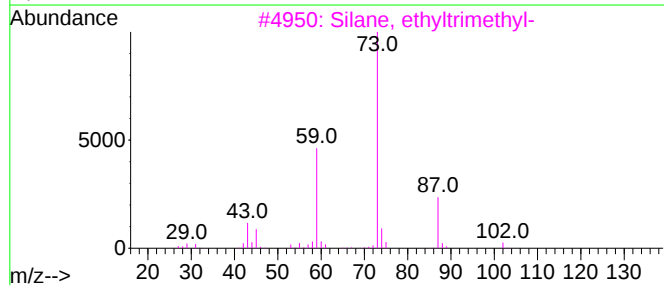
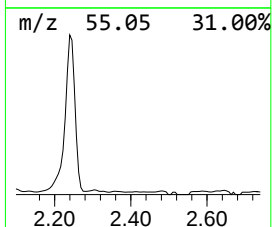
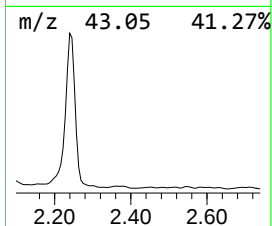
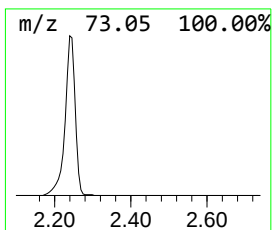
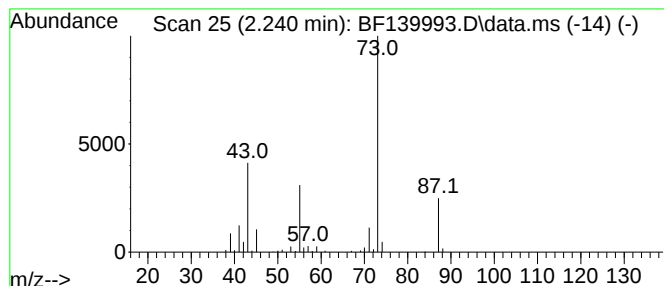
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.240	3.58 ng	160889	1,4-Dichlorobenzene-d4	6.886

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	39
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17



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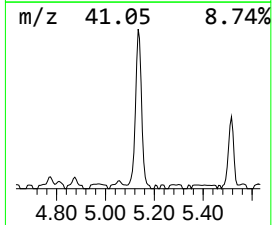
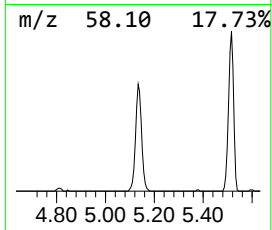
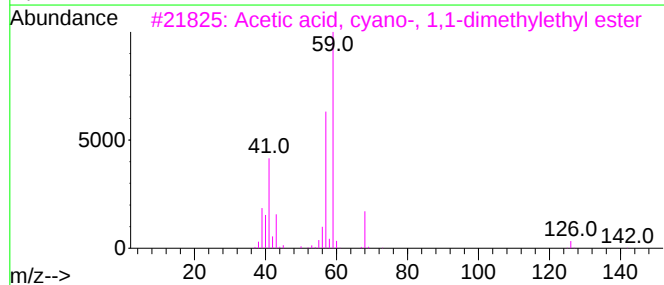
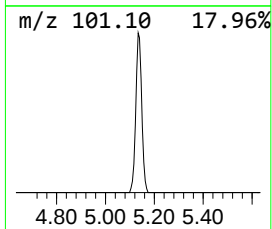
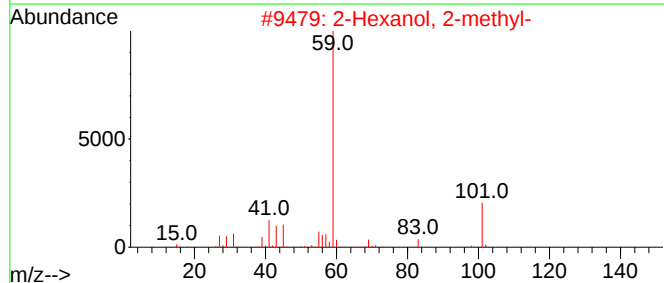
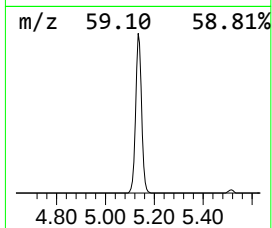
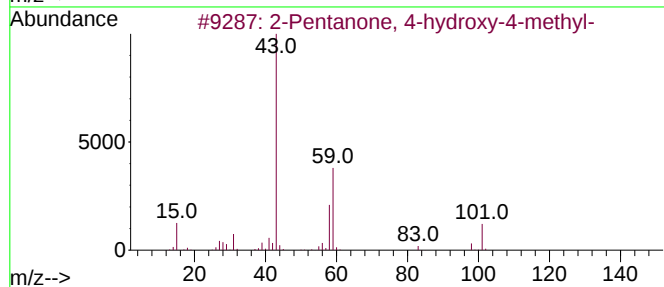
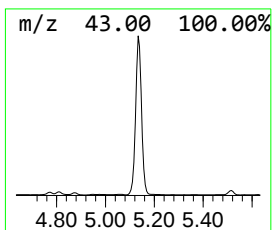
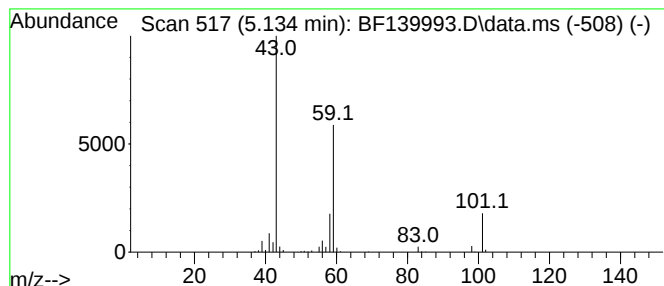
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.134	7.53 ng	338596	1,4-Dichlorobenzene-d4			6.886
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	28
3	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
4	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	12
5	Morpholine, 4-methyl-		101	C5H11NO	000109-02-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.240	3.6	ng	160889	1	6.886	898847	20.0
2-Pentanone, 4-...	5.134	7.5	ng	338596	1	6.886	898847	20.0