

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
 Data File : BF139924.D
 Acq On : 22 Oct 2024 13:01
 Operator : RC/JU
 Sample : PB164246BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164246BL

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: BF139924.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	12	25	31	rBV	76387	138638	2.53%	0.409%
2	5.134	509	517	524	rBV	198585	336613	6.13%	0.993%
3	5.516	574	582	587	rBV	3175579	4652539	84.78%	13.729%
4	6.510	743	751	756	rBV	3108673	4667203	85.05%	13.772%
5	6.887	810	815	821	rVB	711311	910837	16.60%	2.688%
6	7.451	904	911	916	rBV	2141866	3107731	56.63%	9.170%
7	8.169	1027	1033	1038	rBV	945650	1214122	22.13%	3.583%
8	9.245	1209	1216	1221	rBV	3993238	5487557	100.00%	16.193%
9	9.927	1325	1332	1337	rBV	1106024	1406206	25.63%	4.149%
10	10.716	1459	1466	1471	rBV	2517638	3358292	61.20%	9.910%
11	11.410	1579	1584	1590	rBV	1139065	1440282	26.25%	4.250%
12	12.998	1848	1854	1860	rBV	3921270	5407751	98.55%	15.957%
13	14.051	2027	2033	2045	rBV	767662	984991	17.95%	2.907%
14	15.527	2278	2284	2296	rBV	496455	776149	14.14%	2.290%

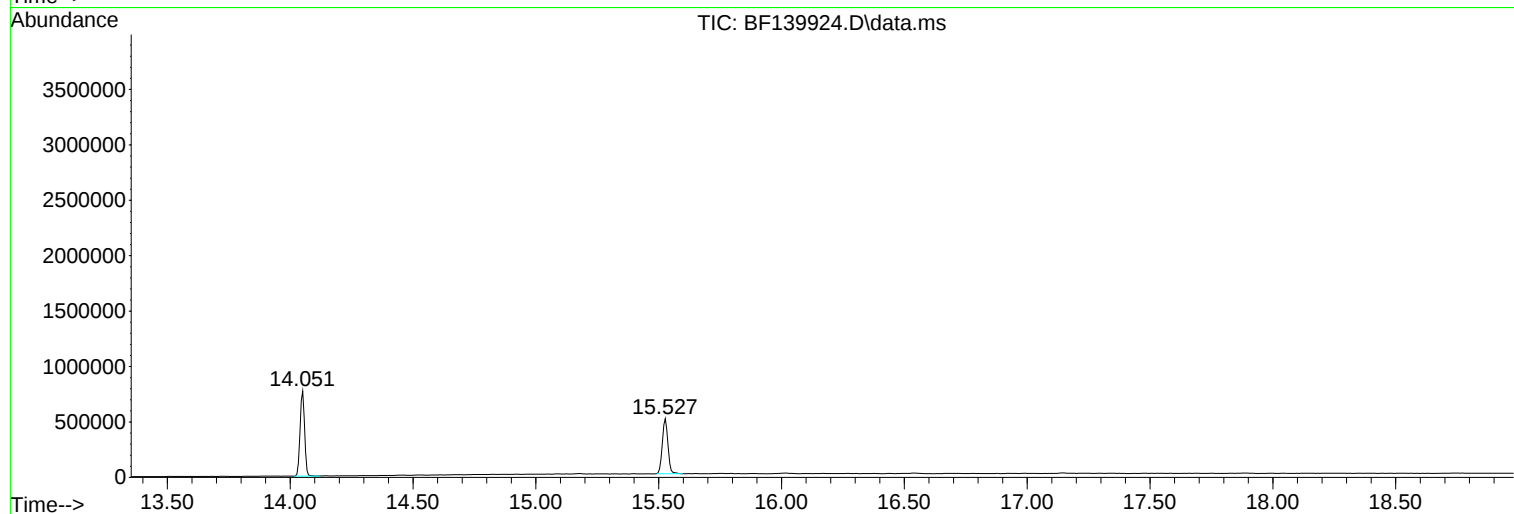
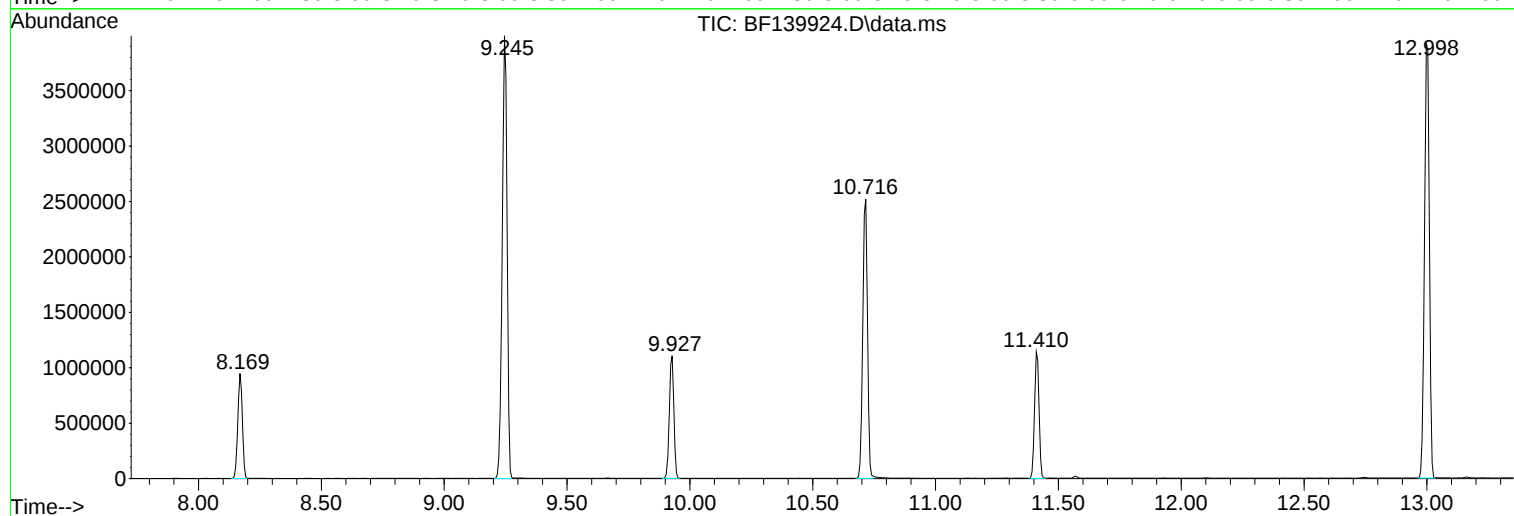
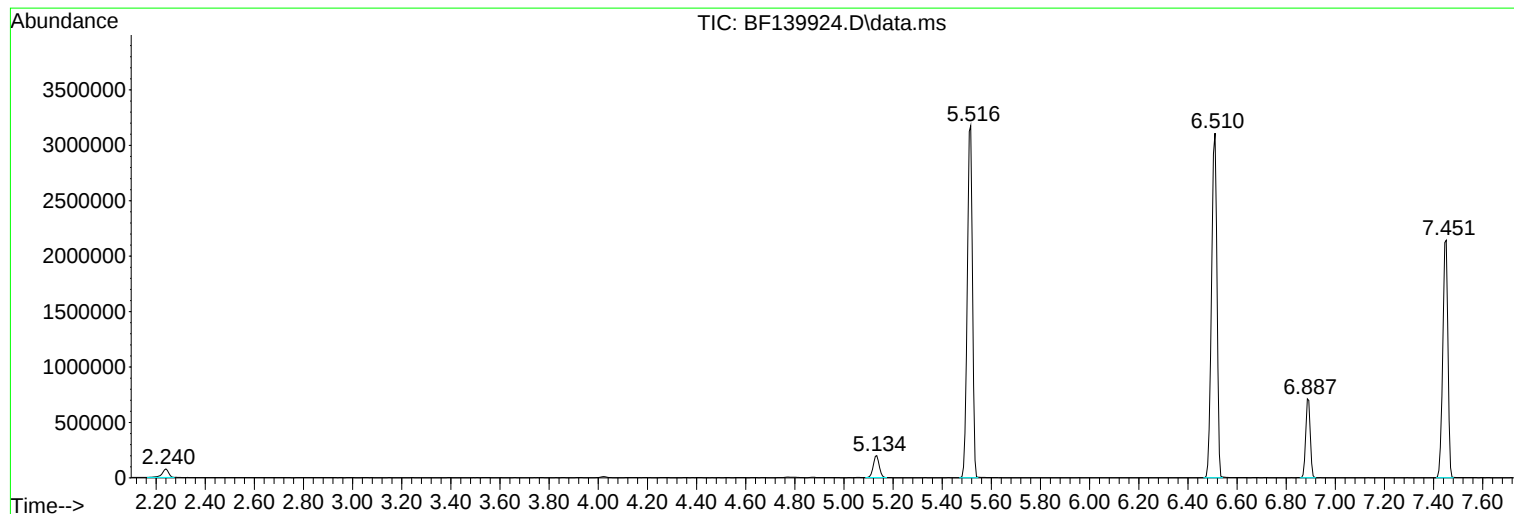
Sum of corrected areas: 33888911

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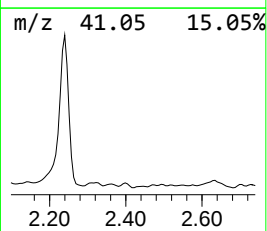
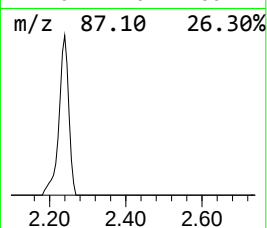
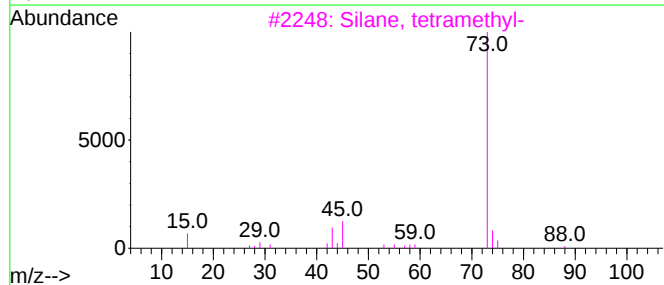
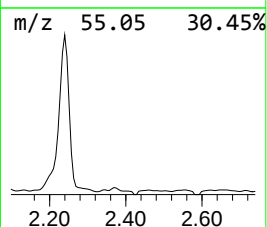
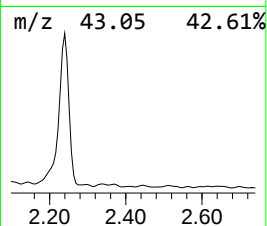
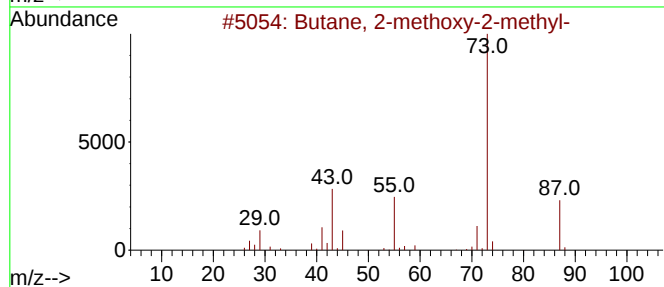
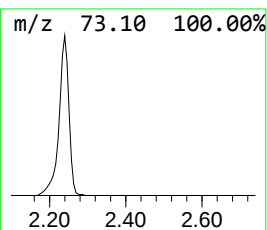
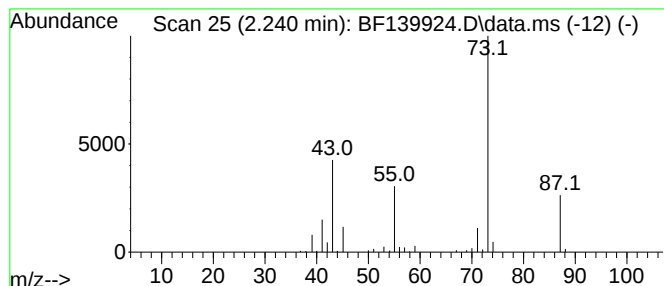
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.240	3.04 ng	138638	1,4-Dichlorobenzene-d4	6.892

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, tetramethyl-	88	C4H12Si	000075-76-3	35
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	12



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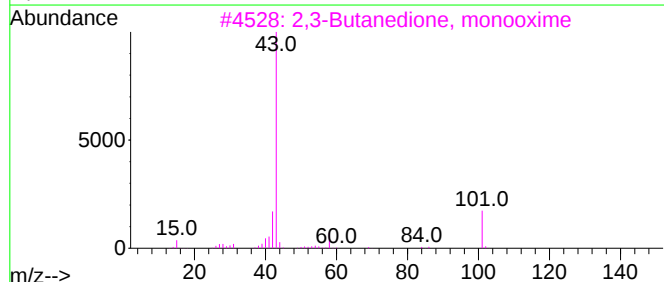
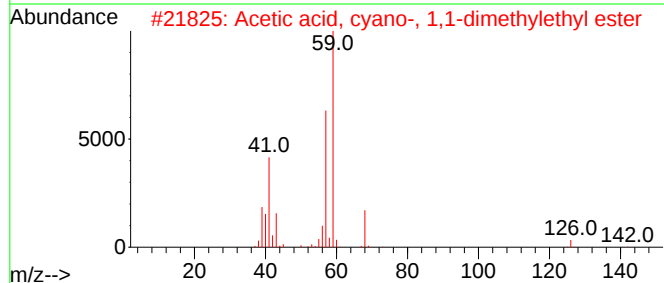
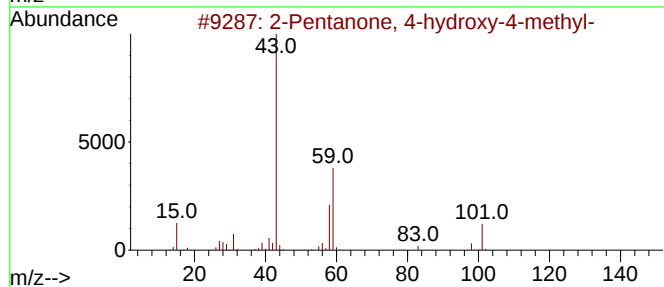
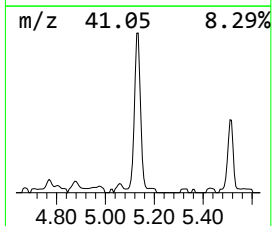
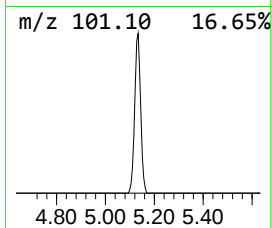
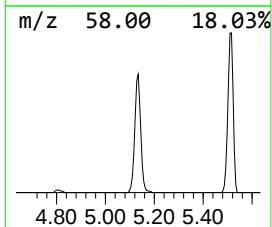
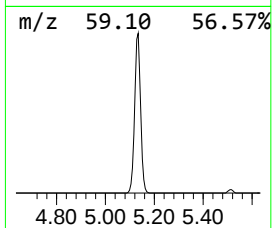
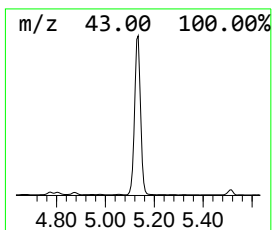
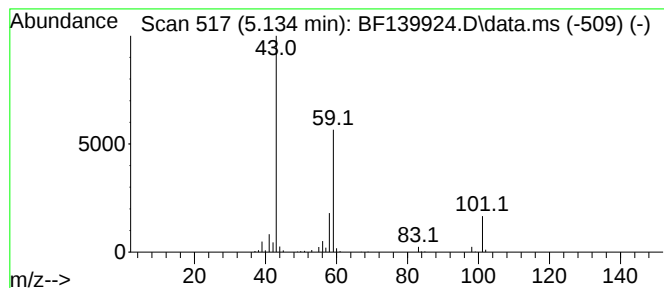
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.134	7.39 ng	336613	1,4-Dichlorobenzene-d4	6.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			N,N'-Bis(2-methyl-2-nitrosopenta...	258	C12H22N2O4	094514-30-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.240	3.0	ng	138638	1	6.892	910837	20.0
2-Pentanone, 4-...	5.134	7.4	ng	336613	1	6.892	910837	20.0