

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
 Data File : BF139962.D
 Acq On : 23 Oct 2024 14:04
 Operator : RC/JU
 Sample : PB164286BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164286BL

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: BF139962.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.246	16	26	32	rVB	96327	169159	3.05%	0.506%
2	5.134	509	517	526	rBV	191313	325538	5.87%	0.974%
3	5.516	575	582	587	rBV	3167895	4525109	81.53%	13.534%
4	6.510	744	751	756	rBV	3045622	4503924	81.15%	13.471%
5	6.887	810	815	821	rVB	704321	884368	15.93%	2.645%
6	7.445	904	910	916	rBV	2114373	3005131	54.15%	8.988%
7	8.169	1027	1033	1038	rBV	919728	1164539	20.98%	3.483%
8	9.245	1209	1216	1221	rBV	3958388	5341336	96.24%	15.975%
9	9.922	1325	1331	1344	rVB	1028668	1359976	24.50%	4.068%
10	10.716	1459	1466	1492	rBV	2440568	3383675	60.97%	10.120%
11	11.410	1579	1584	1590	rBV	1126626	1432080	25.80%	4.283%
12	12.998	1848	1854	1860	rBV	4031941	5550060	100.00%	16.599%
13	14.051	2027	2033	2044	rBV	751869	980332	17.66%	2.932%
14	15.527	2278	2284	2299	rBV	505903	809910	14.59%	2.422%

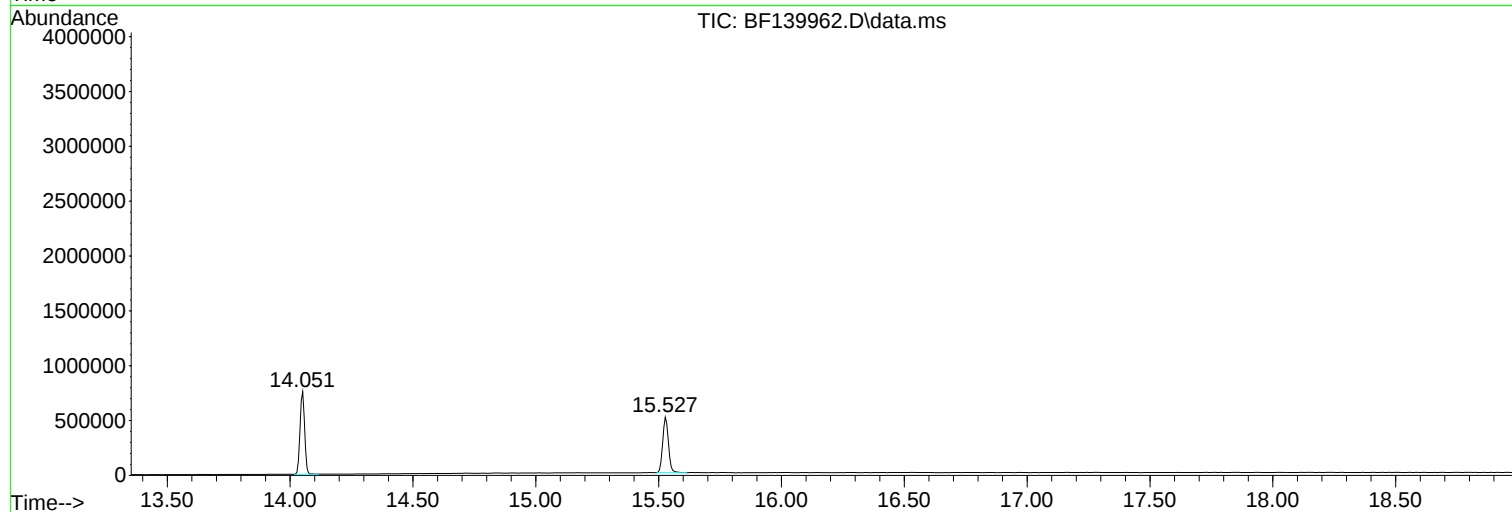
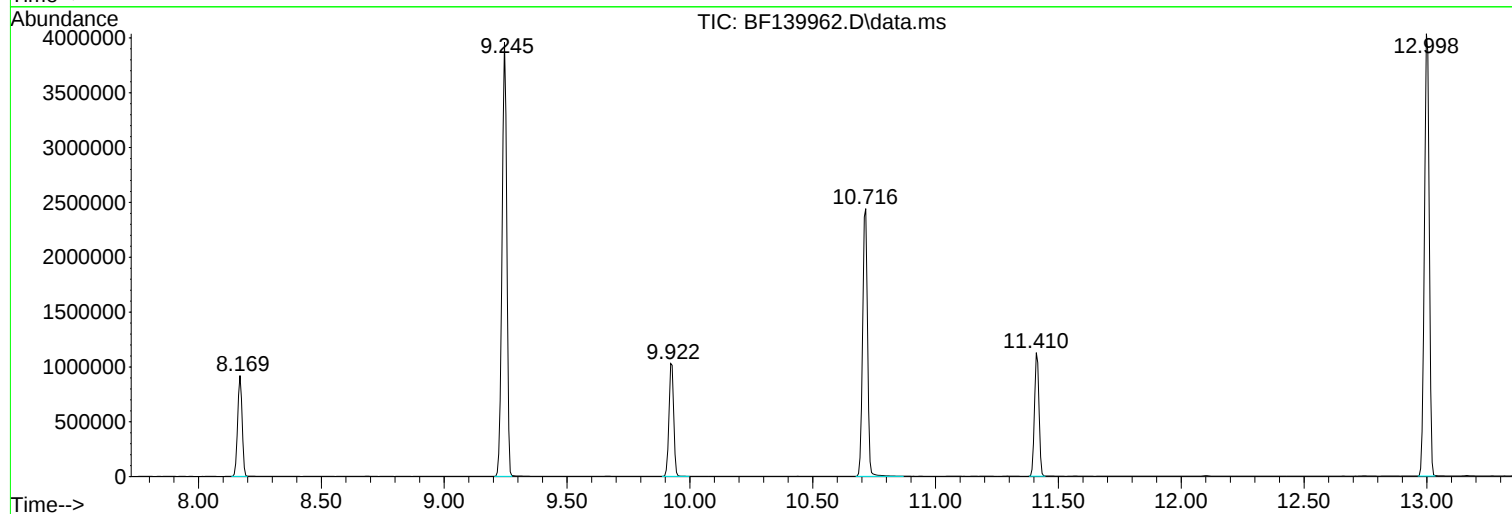
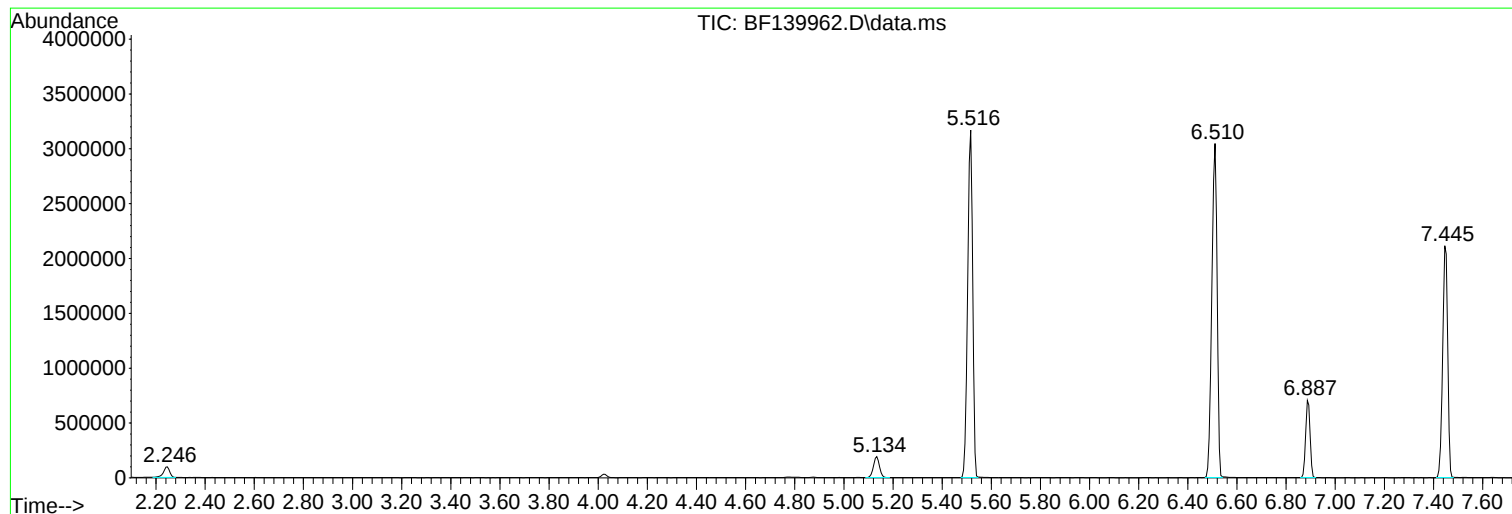
Sum of corrected areas: 33435137

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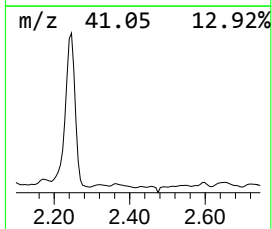
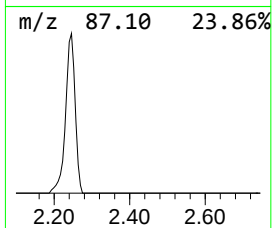
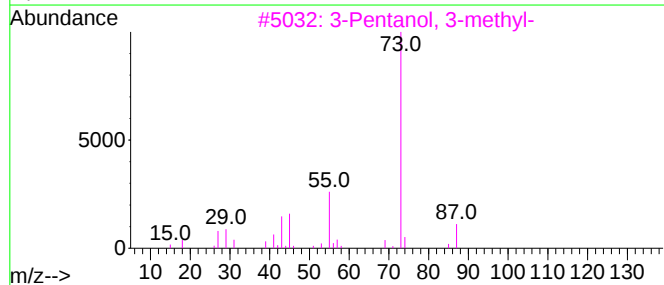
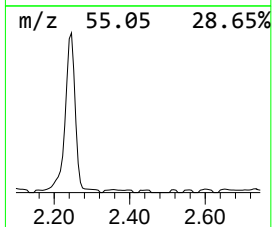
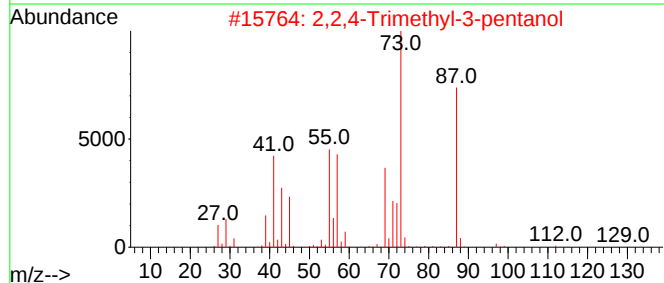
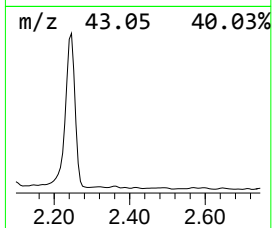
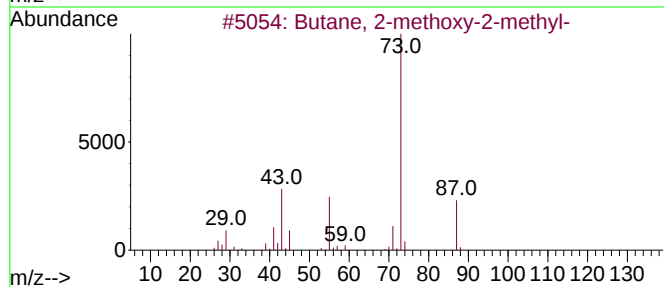
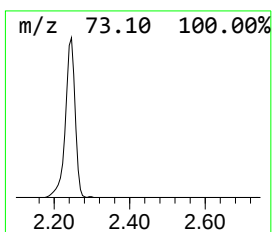
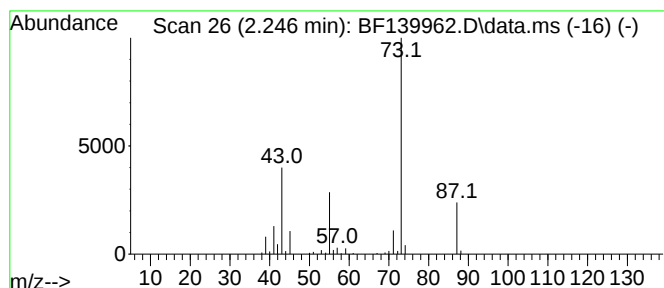
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.246	3.83 ng	169159	1,4-Dichlorobenzene-d4	6.887

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			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	25
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	12



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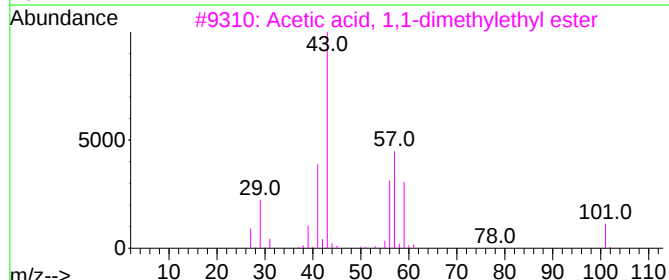
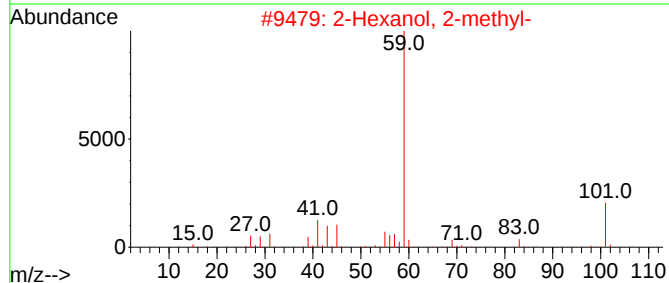
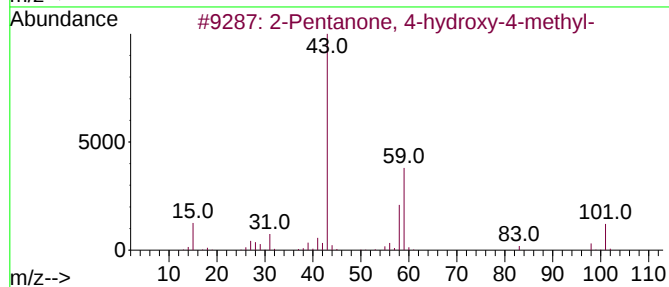
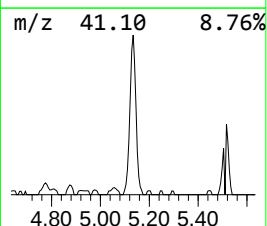
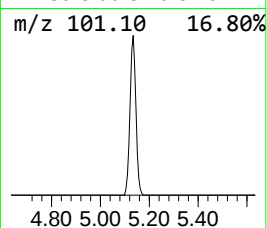
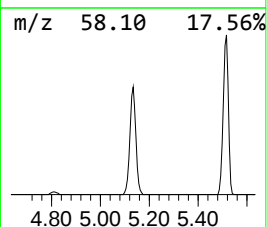
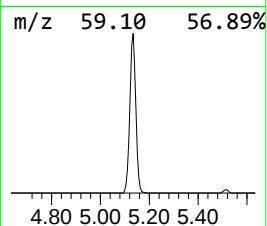
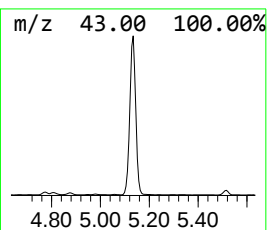
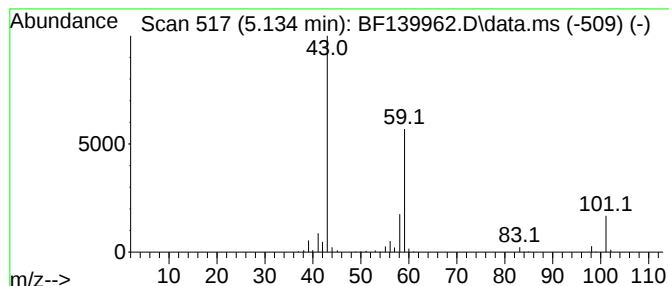
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.134	7.36 ng	325538	1,4-Dichlorobenzene-d4	6.887

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, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.246	3.8	ng	169159	1	6.887	884368	20.0
2-Pentanone, 4-...	5.134	7.4	ng	325538	1	6.887	884368	20.0