

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121322\
 Data File : BF131611.D
 Acq On : 13 Dec 2022 14:18
 Operator : CG\JU
 Sample : PB149584BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB149584BL

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-BF120822.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131611.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.193	10	18	24	rVB	48642	63126	2.36%	0.451%
2	5.104	507	513	522	rVB	441153	529388	19.76%	3.784%
3	5.504	577	581	585	rBV	1428165	1367343	51.03%	9.774%
4	6.516	748	753	756	rBV	1512188	1374911	51.31%	9.828%
5	6.881	811	815	818	rBV	512088	394796	14.73%	2.822%
6	7.451	906	912	915	rBV	1423280	1330411	49.65%	9.510%
7	8.169	1029	1034	1038	rBV	618077	522564	19.50%	3.735%
8	9.251	1212	1218	1221	rBV	2604449	2500011	93.29%	17.870%
9	9.933	1329	1334	1338	rBV	766831	632689	23.61%	4.522%
10	10.728	1465	1469	1472	rBV	1230253	1008975	37.65%	7.212%
11	11.427	1584	1588	1592	rBV	785629	680360	25.39%	4.863%
12	13.027	1855	1860	1863	rBV	3040844	2679705	100.00%	19.154%
13	14.080	2035	2039	2043	rBV	565051	517522	19.31%	3.699%
14	15.592	2290	2296	2302	rBV	295526	388314	14.49%	2.776%

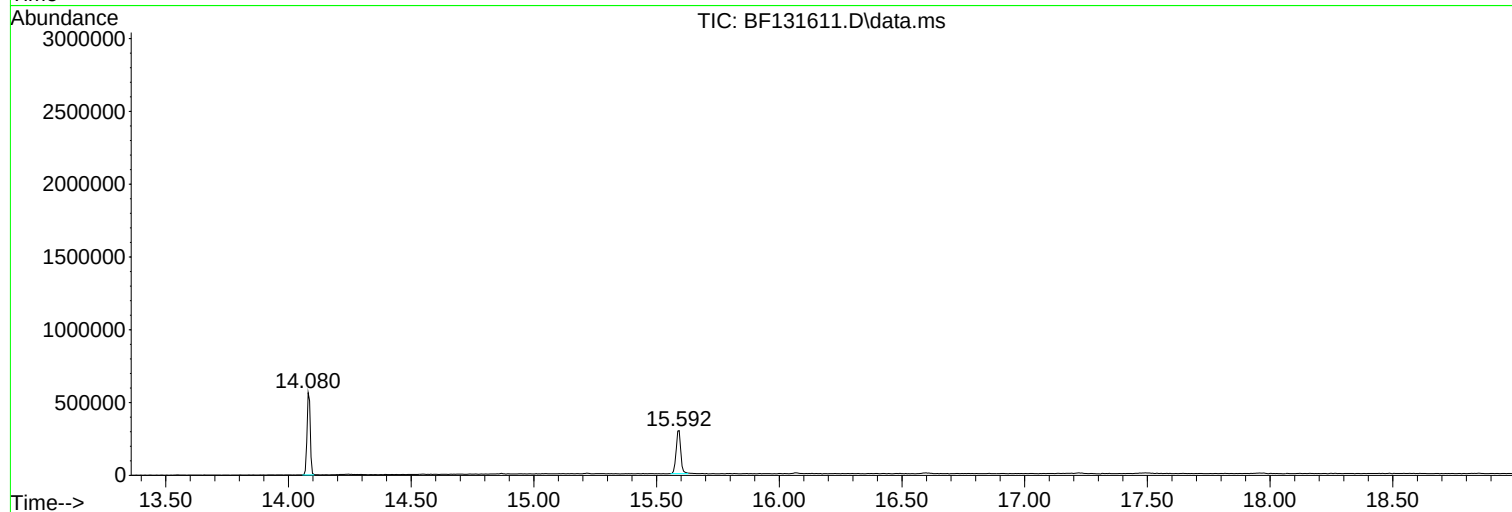
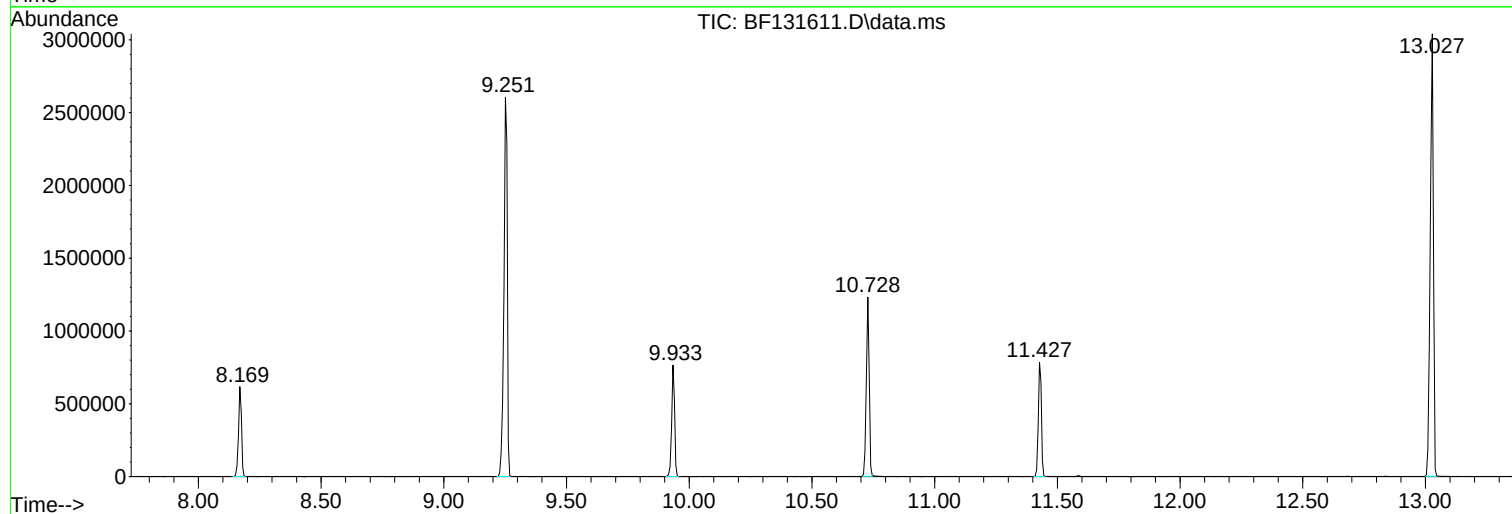
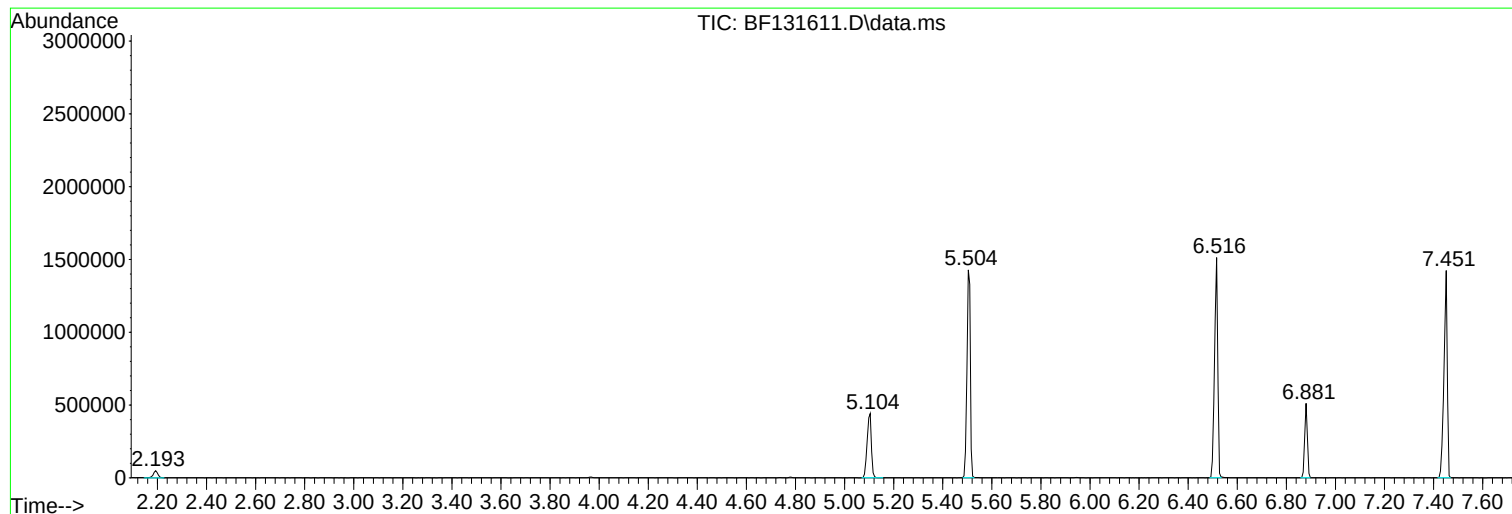
Sum of corrected areas: 13990115

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120822.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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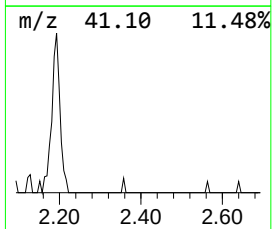
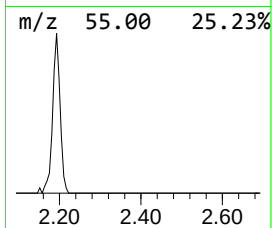
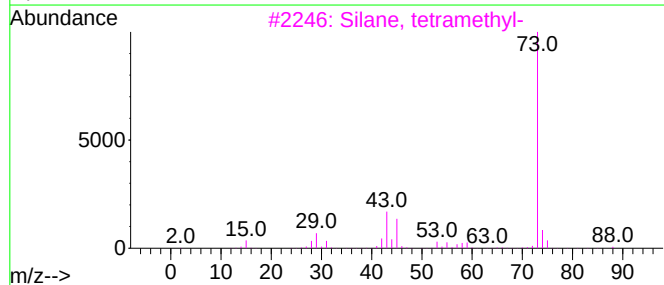
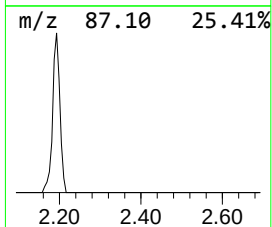
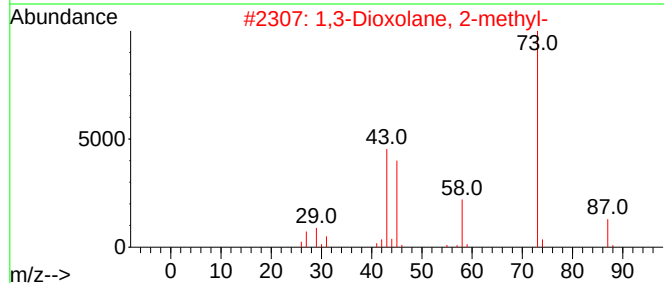
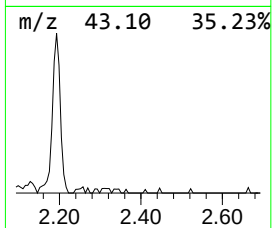
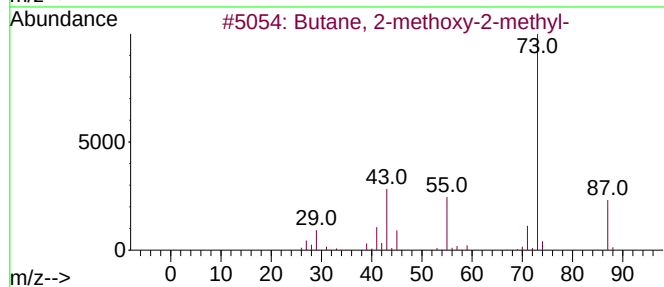
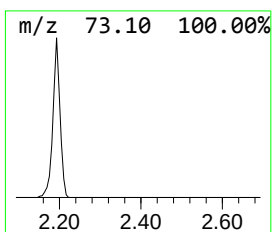
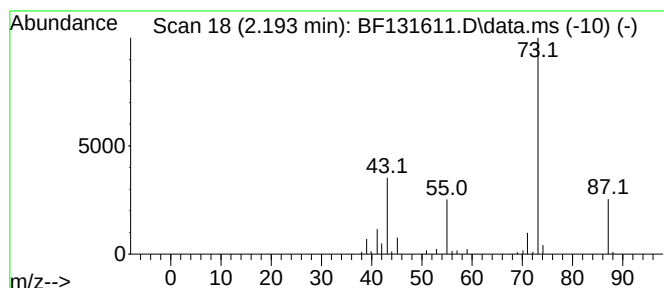
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.193	3.20 ng	63126	1,4-Dichlorobenzene-d4	6.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17



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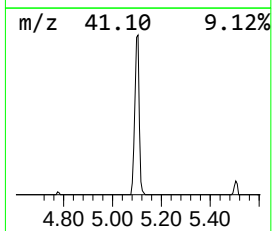
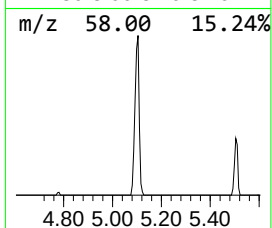
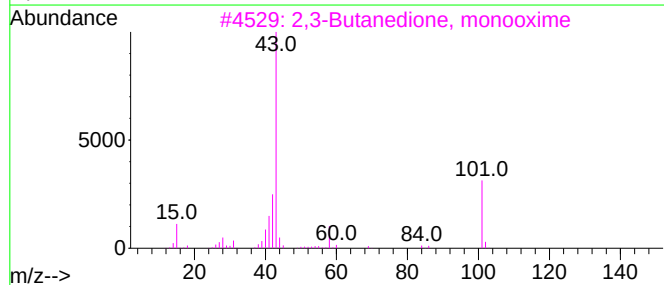
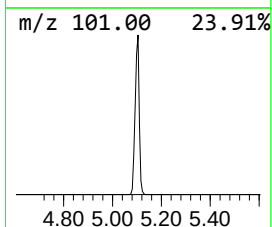
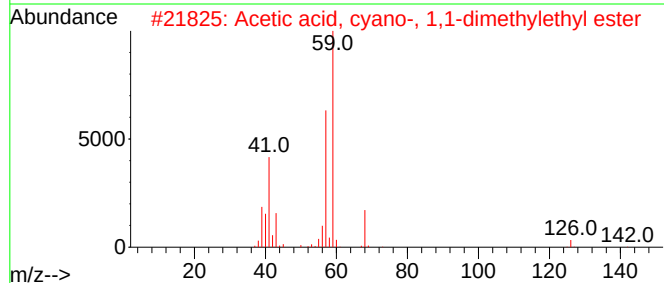
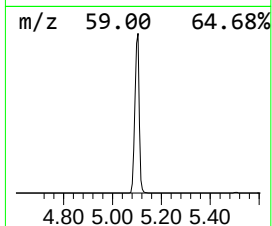
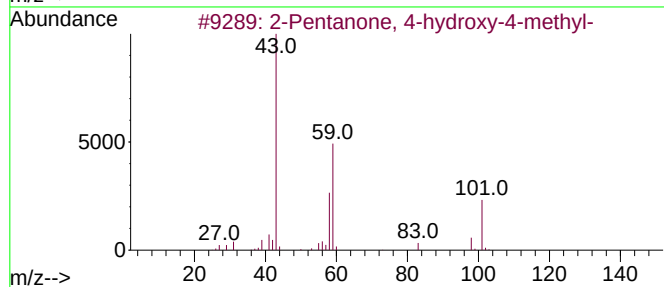
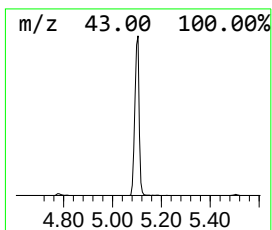
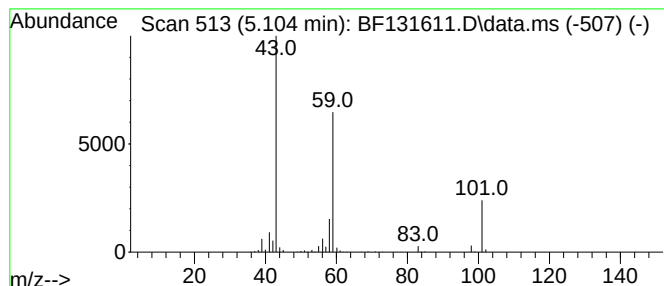
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.104	26.82 ng	529388	1,4-Dichlorobenzene-d4	6.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.193	3.2	ng	63126	1	6.881	394796	20.0
2-Pentanone, 4-...	5.104	26.8	ng	529388	1	6.881	394796	20.0