

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112621\
 Data File : BM033272.D
 Acq On : 26 Nov 2021 19:29
 Operator : CG/JU
 Sample : M4815-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FB

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_M\Methods\8270-BM112421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM033272.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.054	287	292	298	rBV	162391	241138	2.77%	0.707%
2	5.507	363	369	387	rVB	1165931	1764624	20.28%	5.174%
3	6.125	467	474	484	rBV	58409	91738	1.05%	0.269%
4	7.095	629	639	654	rBV	673188	1145434	13.16%	3.359%
5	7.931	774	781	788	rBV	463405	762545	8.76%	2.236%
6	9.095	971	979	993	rBV	1824229	3104521	35.68%	9.103%
7	10.730	1246	1257	1267	rVB	603728	1051418	12.08%	3.083%
8	13.177	1666	1673	1686	rBV	4168686	6511513	74.84%	19.094%
9	14.554	1898	1907	1916	rBV	878798	1355036	15.57%	3.973%
10	16.036	2153	2159	2173	rBV	2978881	4438291	51.01%	13.014%
11	17.289	2366	2372	2379	rBV	1062186	1568153	18.02%	4.598%
12	17.412	2388	2393	2398	rVB4	74813	109409	1.26%	0.321%
13	18.177	2519	2523	2528	rBV4	50011	92567	1.06%	0.271%
14	19.900	2811	2816	2826	rVB	6931299	8700979	100.00%	25.514%
15	21.453	3075	3080	3084	rBV	1208495	1640008	18.85%	4.809%
16	23.783	3471	3476	3483	rBV2	817886	1525803	17.54%	4.474%

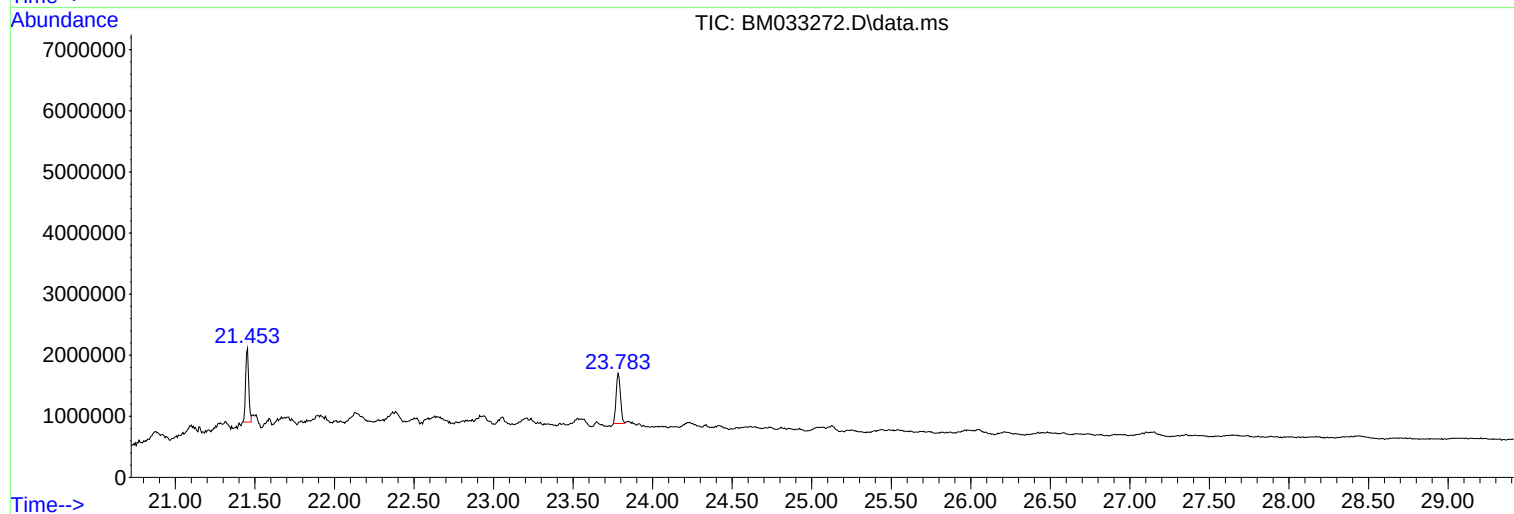
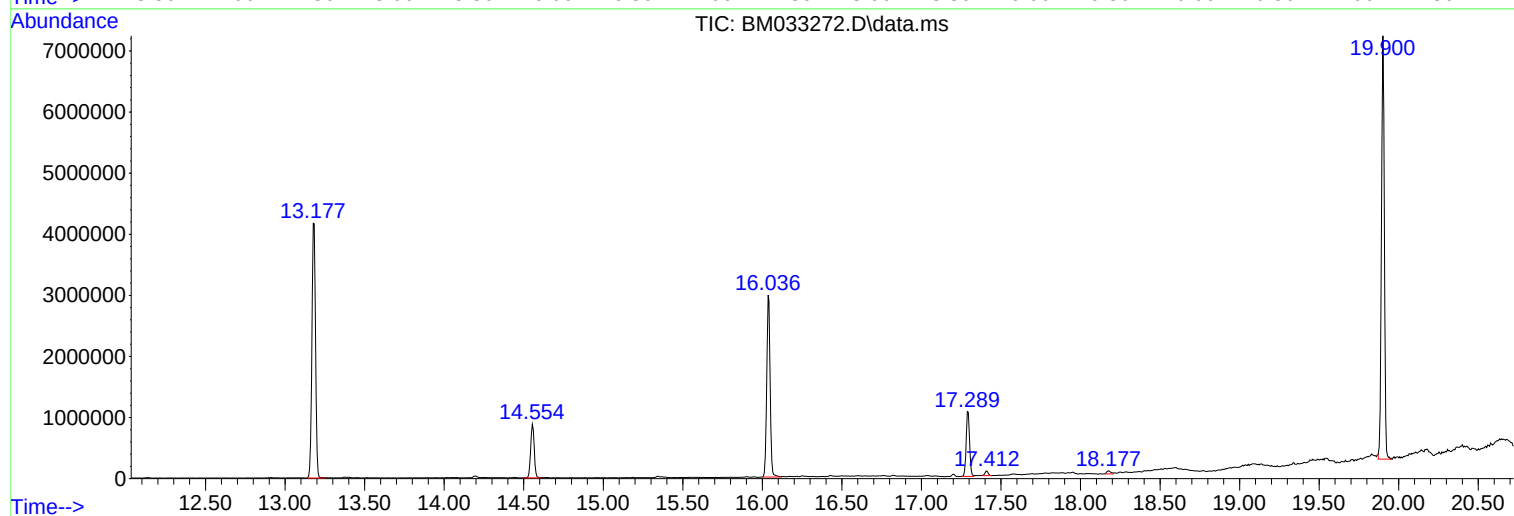
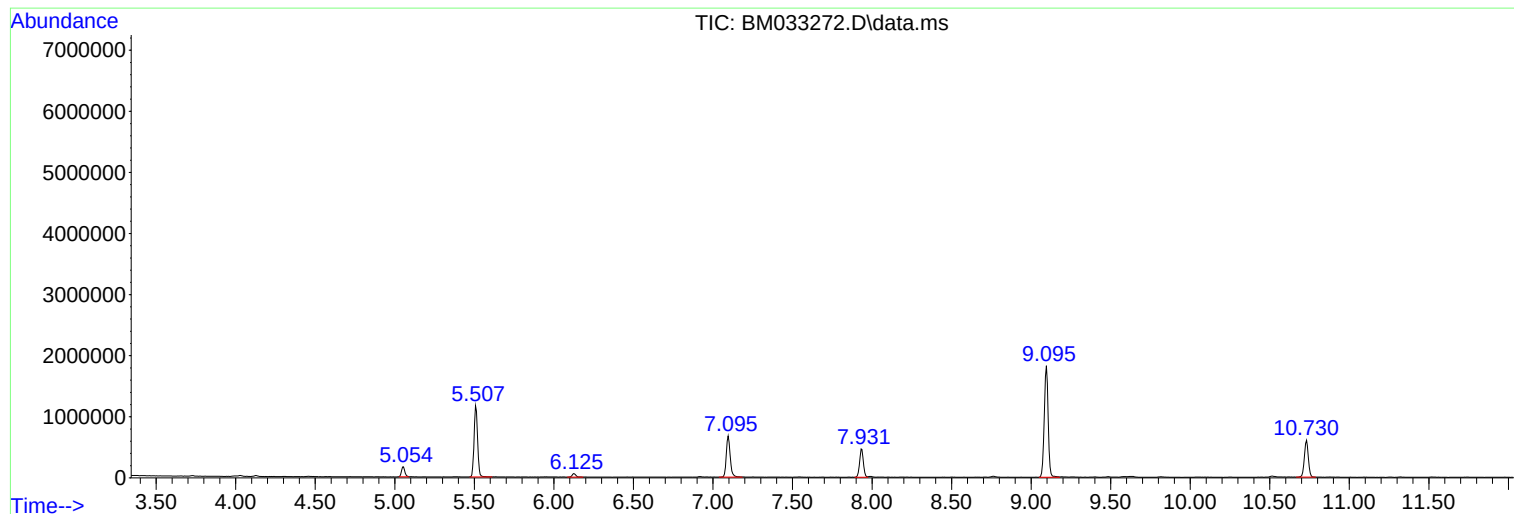
Sum of corrected areas: 34103177

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.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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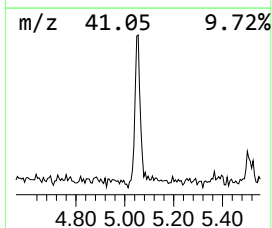
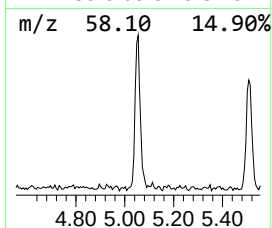
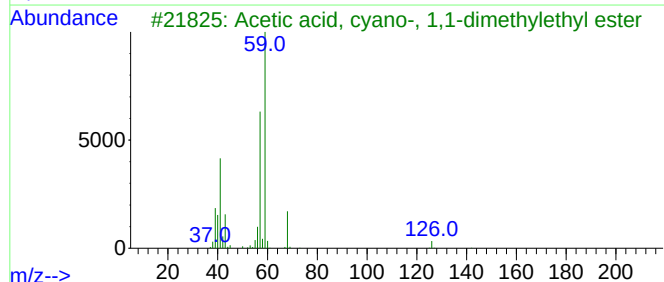
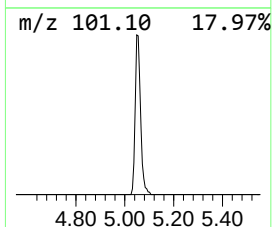
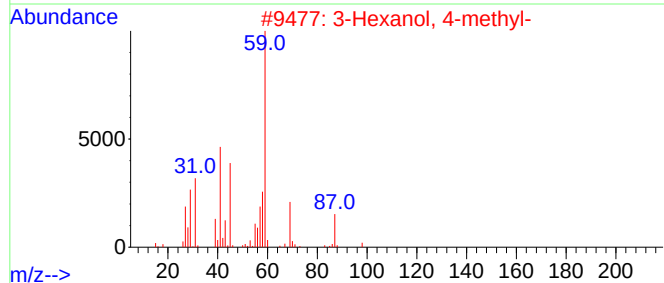
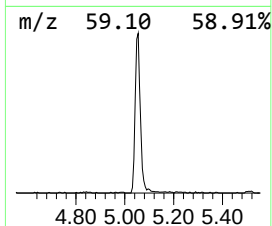
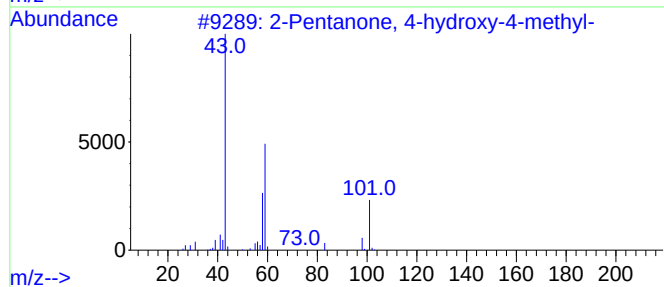
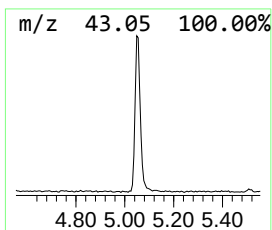
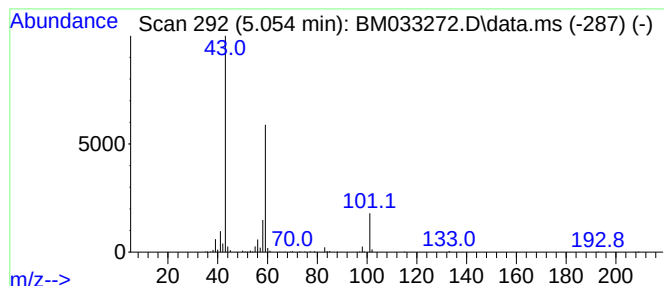
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.054	6.32 ng	241138	1,4-Dichlorobenzene-d4	7.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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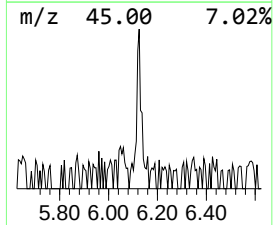
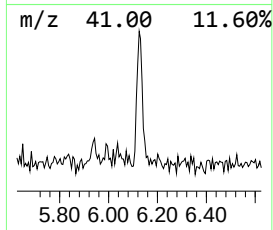
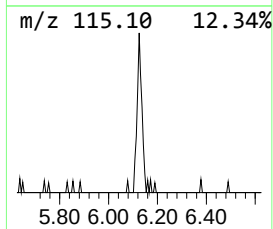
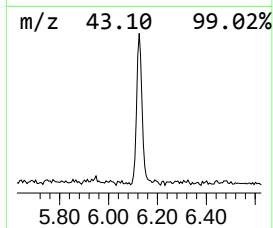
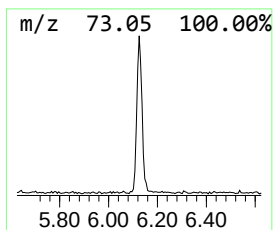
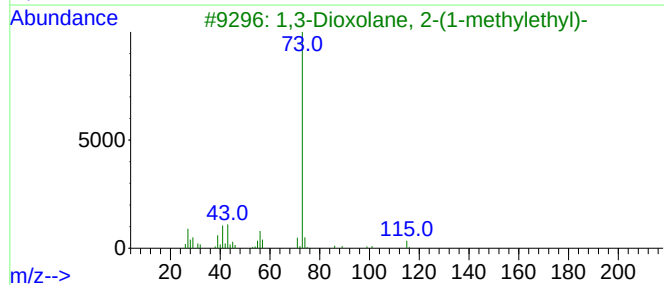
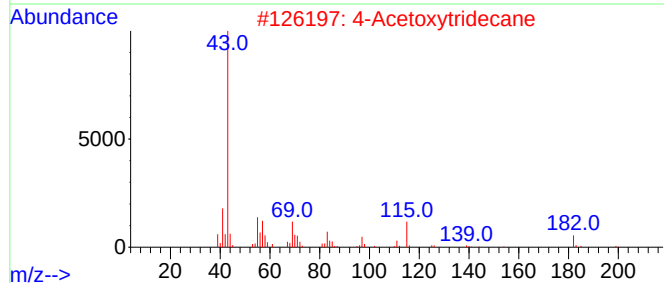
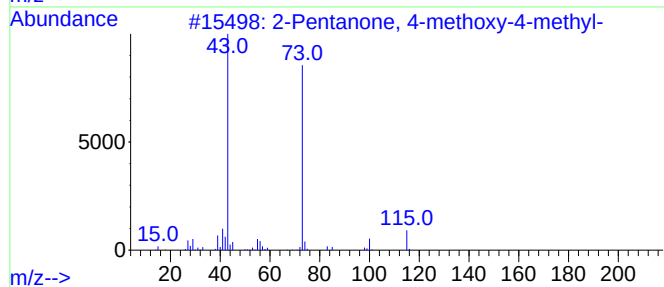
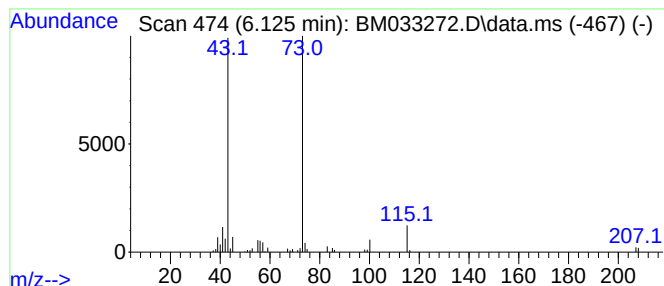
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Peak Number 2 unknown6.125 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.125	2.41 ng	91738	1,4-Dichlorobenzene-d4	7.931

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	72		
2	4-Acetoxytridecane	242	C15H30O2	060826-30-4	23		
3	1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	22		
4	Butane, 2-methoxy-2,3,3-trimethyl-	130	C8H18O	027705-21-1	17		
5	2-Methyl-2,4-dimethoxybutane	132	C7H16O2	039836-89-0	12		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.054	6.3	ng	241138	1	7.931	762545	20.0
unknown6.125	6.125	2.4	ng	91738	1	7.931	762545	20.0