

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082422\
 Data File : BF129983.D
 Acq On : 24 Aug 2022 21:32
 Operator : CG\JU
 Sample : N4262-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-102S

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-BF081422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129983.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.987	456	459	471	rBV	44902	64191	1.72%	0.372%
2	5.404	527	530	553	rBV	1157998	1402843	37.65%	8.130%
3	6.428	701	704	721	rBV	768420	832242	22.34%	4.823%
4	6.763	758	761	770	rBV	705223	626175	16.81%	3.629%
5	7.334	854	858	870	rBV	2098222	2055298	55.16%	11.912%
6	8.045	976	979	993	rVB	1026227	834305	22.39%	4.835%
7	9.122	1157	1162	1165	rBV	4420541	3725957	100.00%	21.594%
8	9.798	1273	1277	1281	rBV	1133938	903291	24.24%	5.235%
9	10.592	1409	1412	1427	rBV	2363145	2126554	57.07%	12.325%
10	11.292	1527	1531	1534	rBV	877763	807575	21.67%	4.680%
11	12.874	1795	1800	1803	rBV	2986826	2595146	69.65%	15.041%
12	13.939	1977	1981	1985	rBV	545818	548114	14.71%	3.177%
13	14.021	1990	1995	2000	rBV	210747	202908	5.45%	1.176%
14	15.386	2223	2227	2234	rVB	459041	529660	14.22%	3.070%

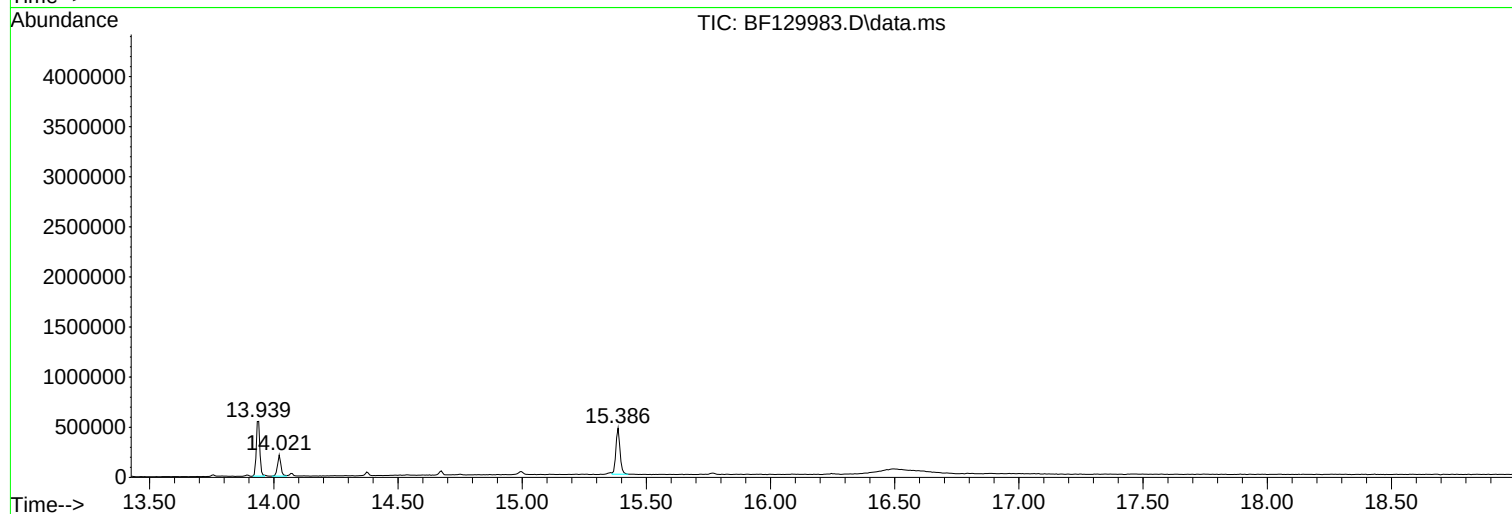
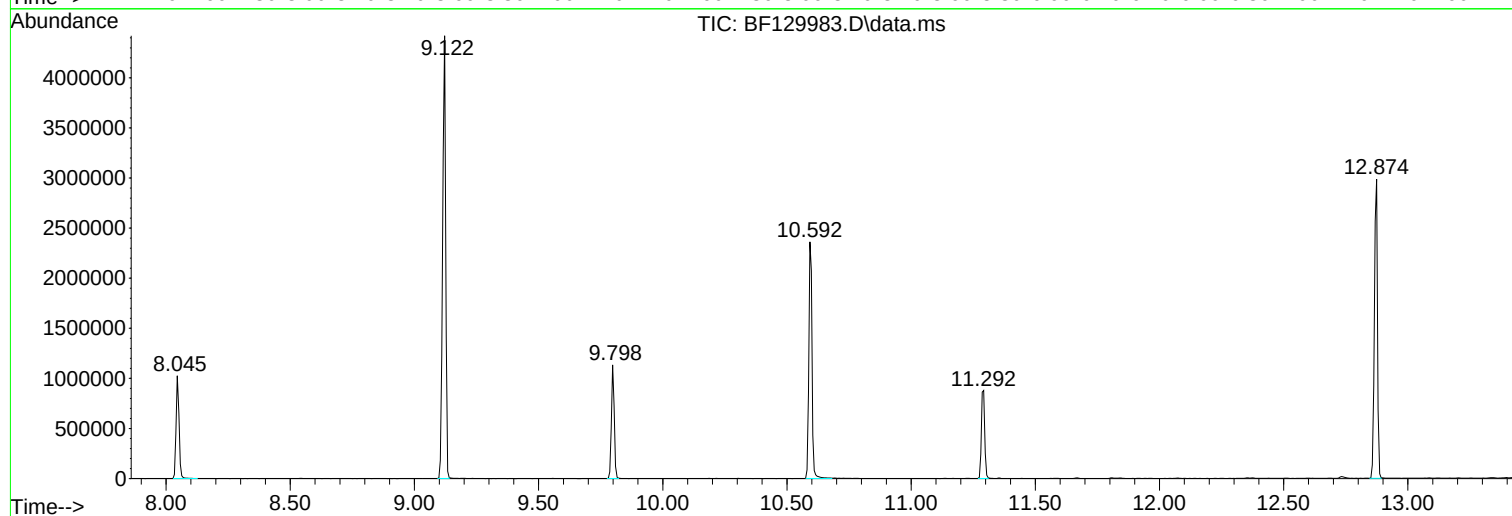
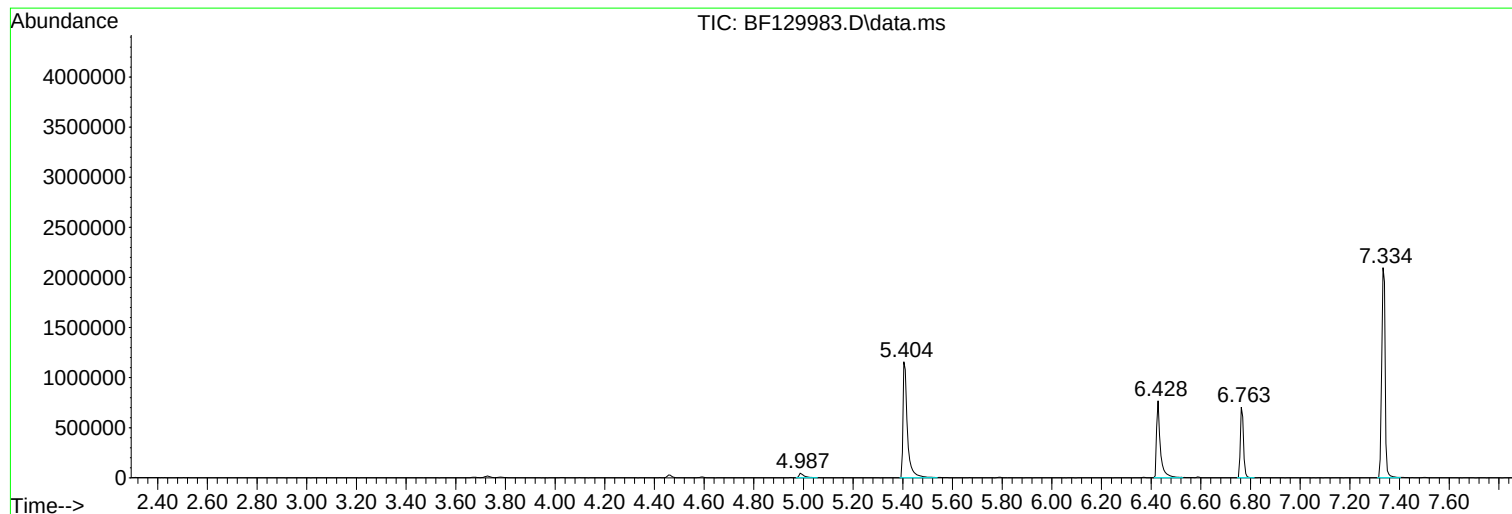
Sum of corrected areas: 17254259

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.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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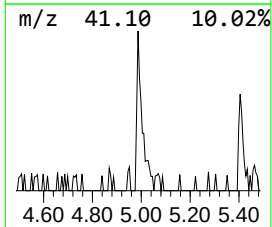
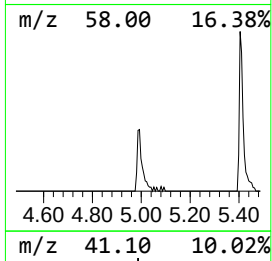
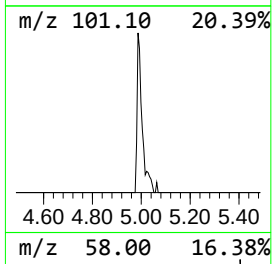
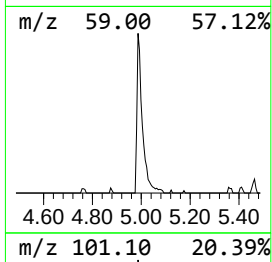
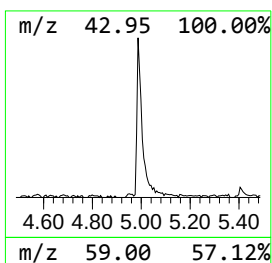
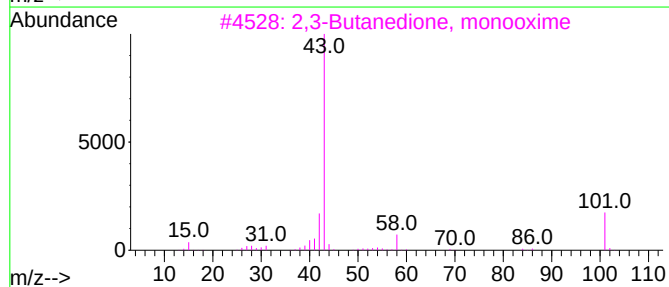
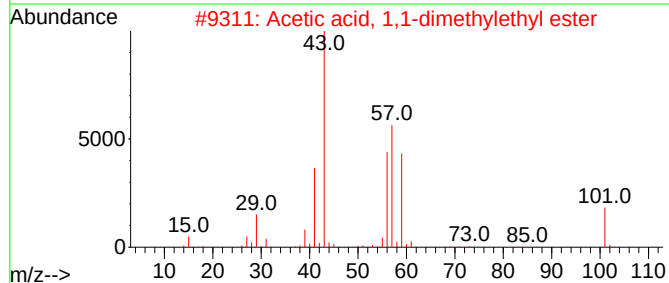
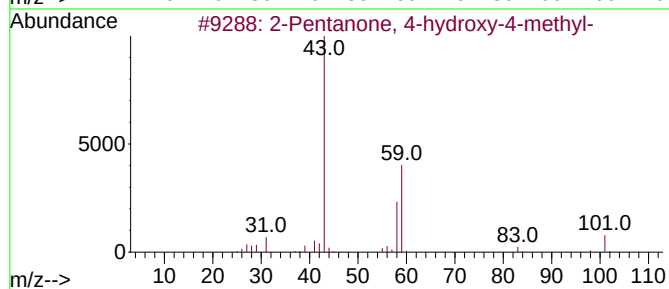
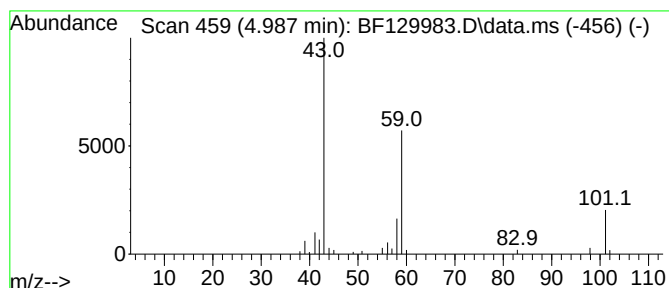
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.987	2.05 ng	64191	1,4-Dichlorobenzene-d4	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
4	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25		
5	1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	17		



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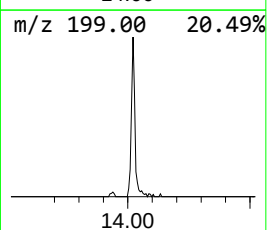
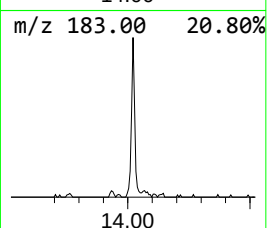
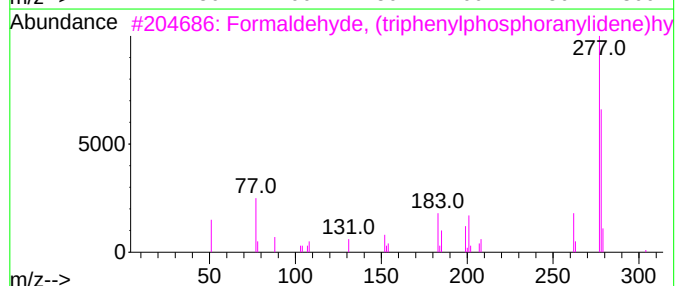
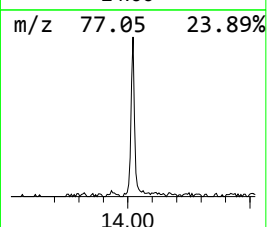
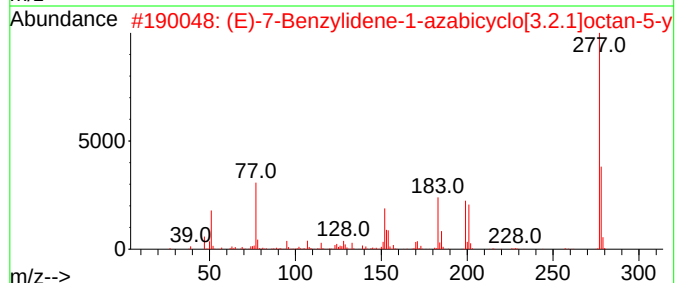
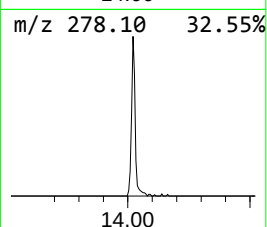
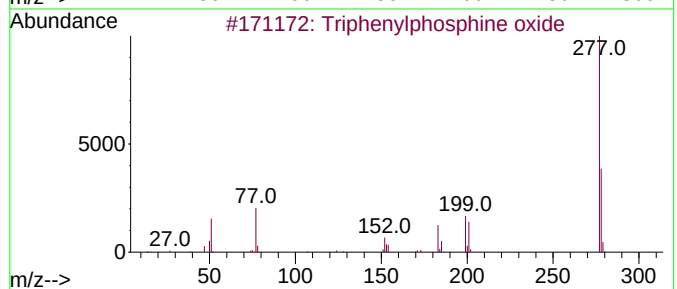
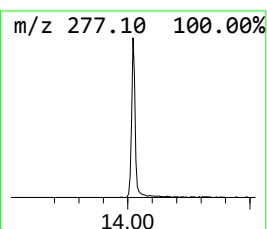
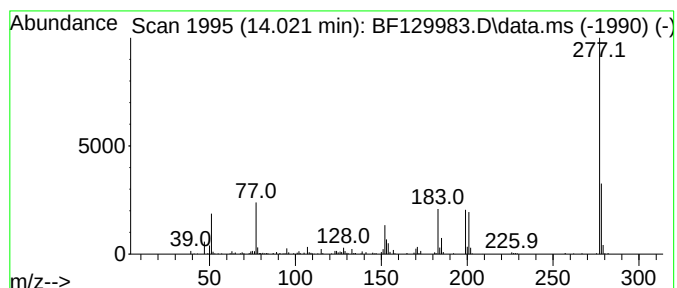
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Peak Number 2 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.021	7.40 ng	202908	Chrysene-d12	13.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	87
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	64
4			Carbazol-1-one, 1,2,3,4-tetrahyd...	277	C18H15N02	299962-12-2	25
5			[1,2,4]Triazolo[1,5-a]pyrimidine...	278	C15H14N6	1000350-69-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.987	2.0	ng	64191	1	6.763	626175	20.0
Triphenylphosph...	14.021	7.4	ng	202908	5	13.939	548114	20.0