

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062221\
 Data File : BP006022.D
 Acq On : 23 Jun 2021 00:46
 Operator : CG/JU
 Sample : M2773-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-6-AOC-2

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_P\Methods\8270E-BP060821.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006022.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.493	402	409	426	rBV	599493	955392	18.23%	5.829%
2	7.093	674	681	706	rBV	523504	933698	17.81%	5.697%
3	7.916	815	821	833	rBV	232373	389334	7.43%	2.376%
4	9.087	1013	1020	1039	rBV	331858	610194	11.64%	3.723%
5	10.446	1243	1251	1262	rBV3	23193	61949	1.18%	0.378%
6	10.722	1289	1298	1315	rBV	297672	518575	9.89%	3.164%
7	13.175	1708	1715	1734	rBV	1045946	1588727	30.31%	9.694%
8	14.545	1941	1948	1966	rBV	435635	690731	13.18%	4.215%
9	16.034	2194	2201	2223	rBV	1059150	1741811	33.23%	10.628%
10	17.122	2376	2386	2390	rBV3	37987	70023	1.34%	0.427%
11	17.292	2407	2415	2428	rBV	541383	860029	16.41%	5.247%
12	18.175	2560	2565	2570	rBV2	35295	52768	1.01%	0.322%
13	19.839	2843	2848	2861	rBV	4327057	5242105	100.00%	31.985%
14	21.075	3054	3058	3067	rBV2	107559	220720	4.21%	1.347%
15	21.363	3102	3107	3121	rBV	864254	1181630	22.54%	7.210%
16	23.769	3510	3516	3530	rVB2	601349	1271688	24.26%	7.759%

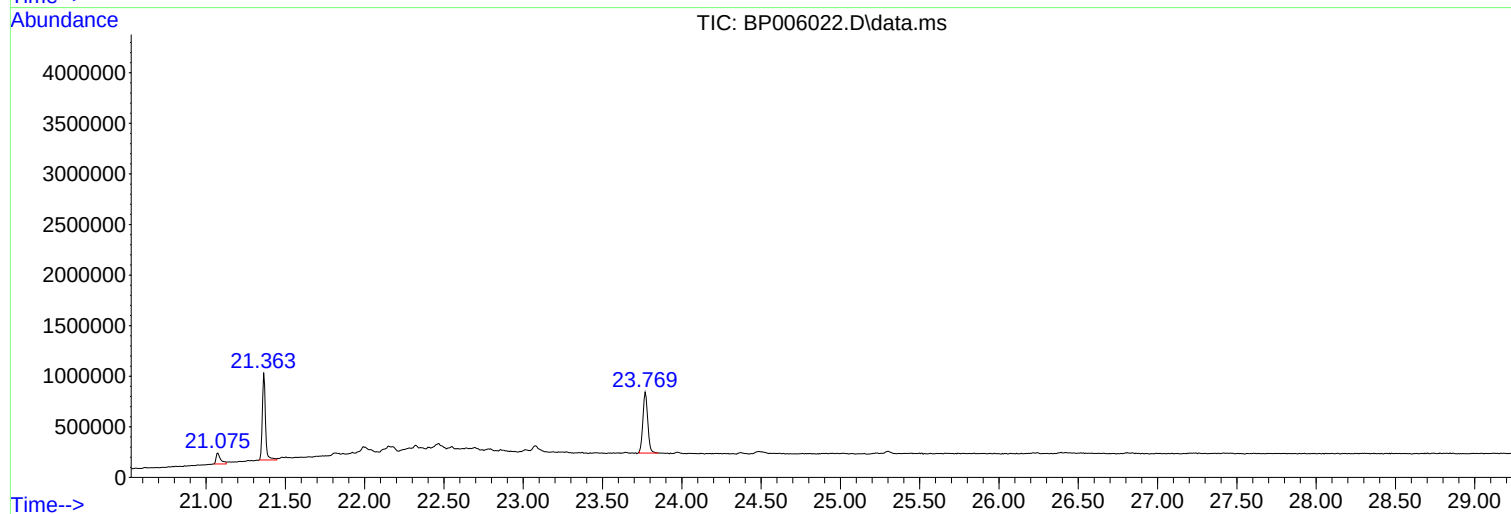
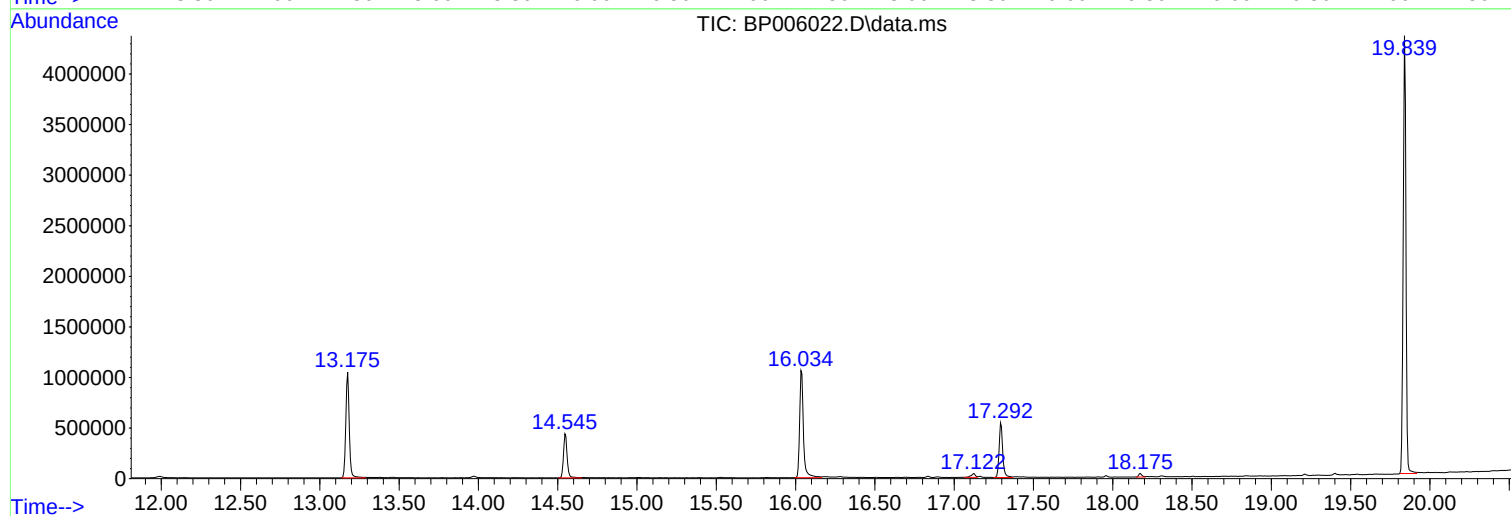
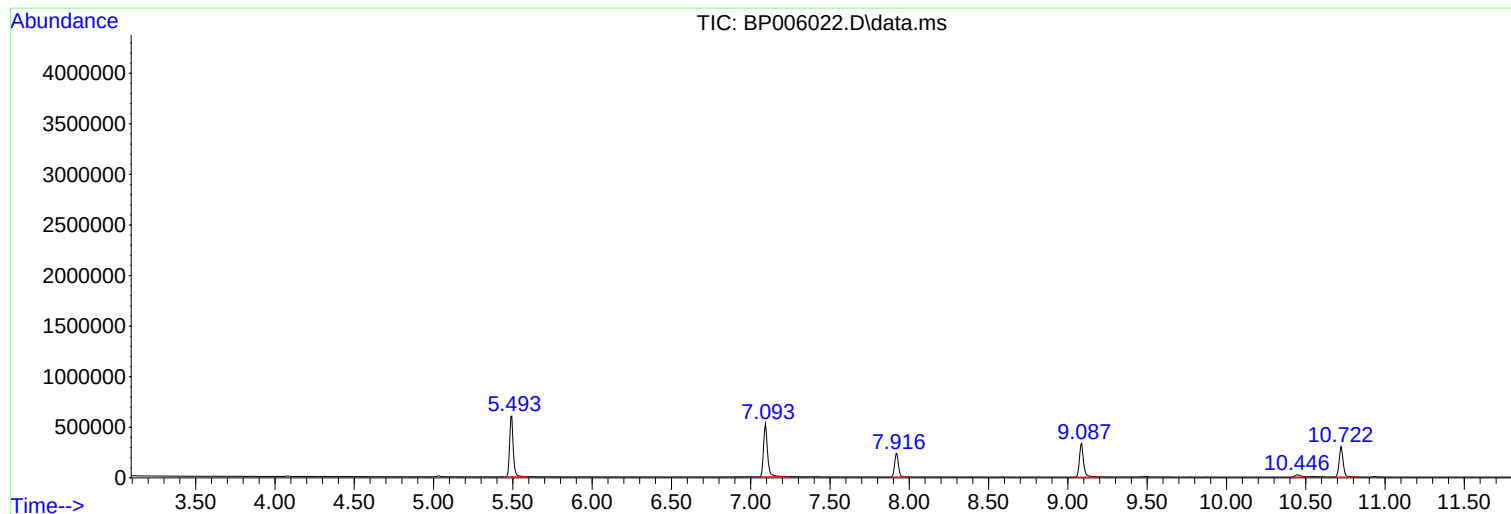
Sum of corrected areas: 16389374

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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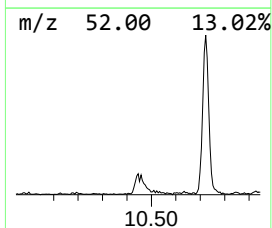
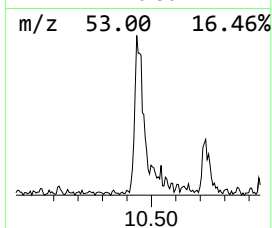
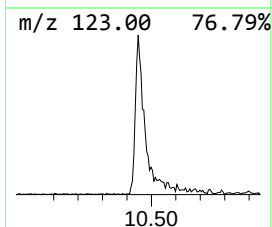
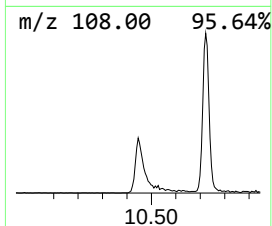
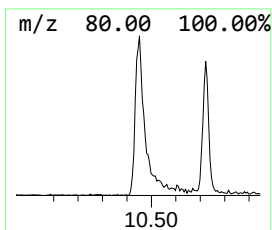
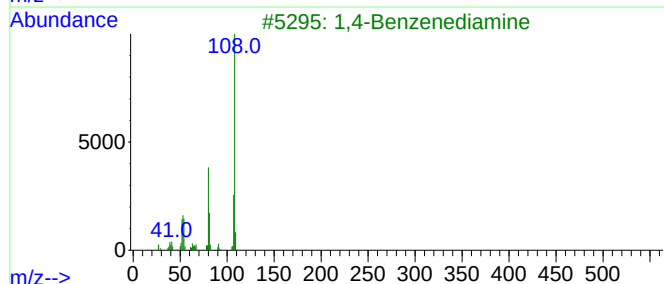
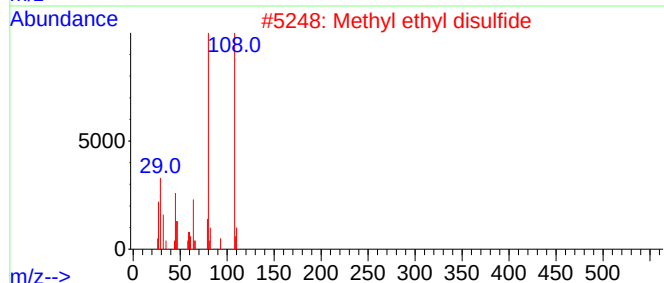
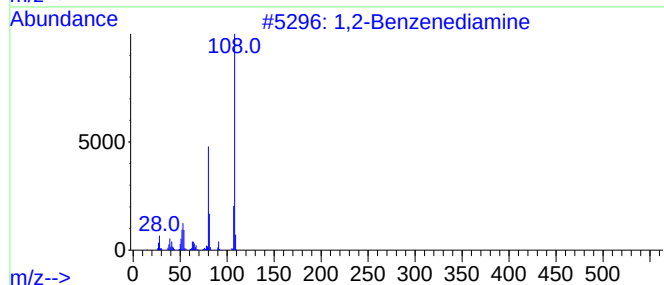
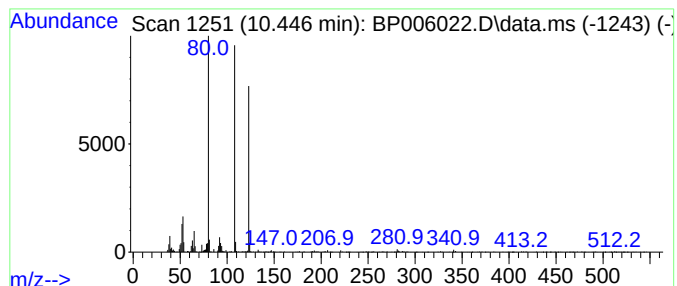
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,2-Benzenediamine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.446	2.39 ng	61949	Naphthalene-d8	10.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenediamine	108	C6H8N2	000095-54-5	72
2			Methyl ethyl disulfide	108	C3H8S2	020333-39-5	64
3			1,4-Benzenediamine	108	C6H8N2	000106-50-3	64
4			3-Pyridinamine, 2-methyl-	108	C6H8N2	003430-10-2	64
5			Tris(ethanesulfonyl)methane	292	C7H16O6S3	021467-59-4	50



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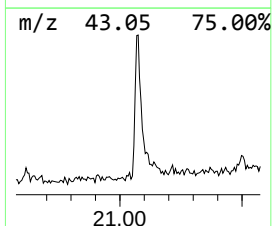
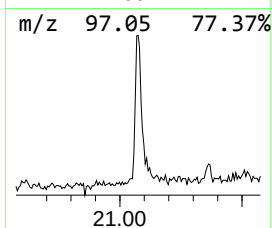
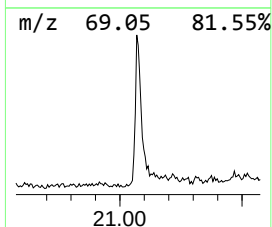
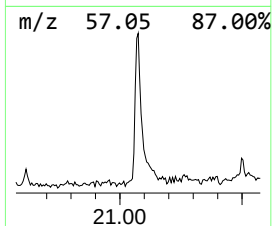
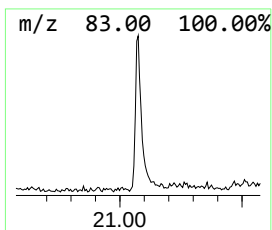
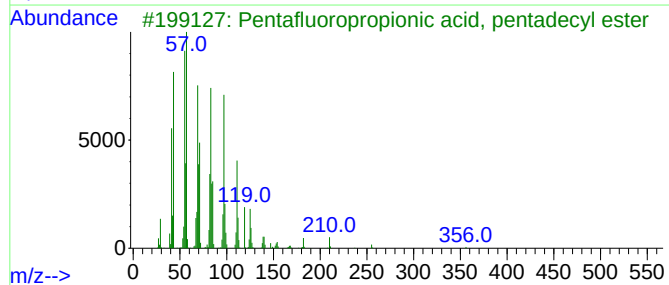
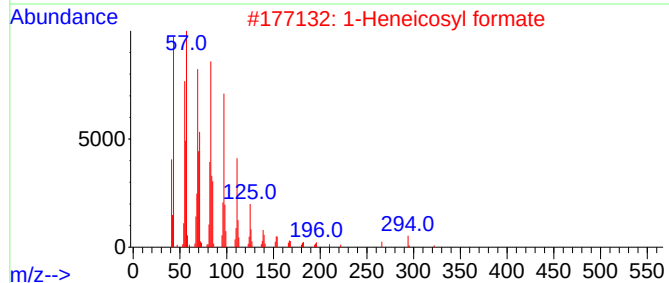
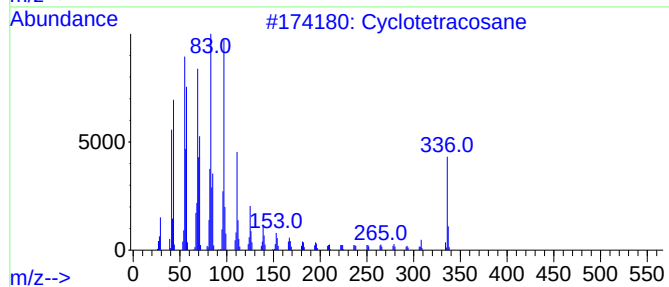
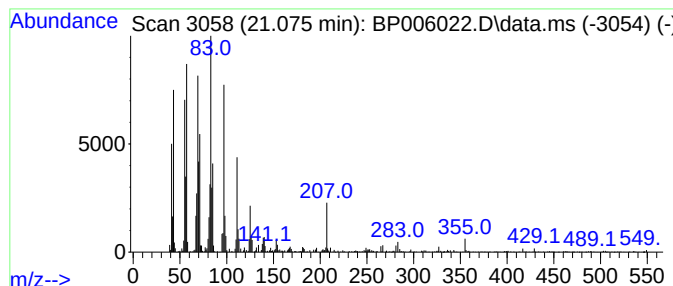
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Peak Number 2 Cyclotetracosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.075	3.74 ng	220720	Chrysene-d12	21.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	94
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	90
3			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	81
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	81
5			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	81



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,2-Benzenediamine	10.446	2.4	ng	61949	2	10.722	518575	20.0
Cyclotetrasiloxane	21.075	3.7	ng	220720	5	21.363	1181630	20.0