

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030722\
 Data File : BF127233.D
 Acq On : 07 Mar 2022 21:21
 Operator : CG\JU
 Sample : N1730-14
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 27-28-SAMPLE-LOCATION-14

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-BF021622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127233.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.504	542	547	550	rBV	2646556	2452172	87.17%	11.702%
2	6.510	713	718	721	rBV	2580094	2546470	90.53%	12.152%
3	6.875	777	780	784	rBV	820370	658009	23.39%	3.140%
4	7.439	871	876	879	rBV	1709994	1724383	61.30%	8.229%
5	7.534	889	892	901	rVB2	24284	32514	1.16%	0.155%
6	8.057	978	981	991	rBV2	17150	41898	1.49%	0.200%
7	8.157	994	998	1002	rBV	991888	909948	32.35%	4.342%
8	9.233	1176	1181	1185	rBV	3083354	2812944	100.00%	13.423%
9	9.304	1188	1193	1197	rVB	42607	37105	1.32%	0.177%
10	9.628	1244	1248	1253	rBV	377947	317185	11.28%	1.514%
11	9.916	1294	1297	1301	rVB	1137482	998918	35.51%	4.767%
12	10.339	1366	1369	1372	rVB2	37111	34826	1.24%	0.166%
13	10.710	1428	1432	1436	rBV	2143867	2096351	74.53%	10.004%
14	10.810	1447	1449	1454	rVB2	31357	37443	1.33%	0.179%
15	11.257	1522	1525	1528	rBV2	36405	38812	1.38%	0.185%
16	11.286	1528	1530	1537	rVB4	23214	37192	1.32%	0.177%
17	11.410	1547	1551	1554	rBV	1302019	1094392	38.91%	5.222%
18	11.433	1554	1555	1558	rVV	189479	113252	4.03%	0.540%
19	11.486	1559	1564	1567	rVB	56822	59237	2.11%	0.283%
20	11.816	1617	1620	1623	rVB	89137	77969	2.77%	0.372%
21	12.627	1754	1758	1761	rBV	195890	172941	6.15%	0.825%
22	12.857	1794	1797	1800	rVB	152938	125631	4.47%	0.600%
23	12.998	1816	1821	1824	rBV	3082787	2719622	96.68%	12.978%
24	13.845	1961	1965	1968	rBV	31983	43124	1.53%	0.206%
25	13.880	1968	1971	1974	rBV	40031	41422	1.47%	0.198%
26	13.921	1974	1978	1981	rBV	108799	108512	3.86%	0.518%
27	13.974	1982	1987	1988	rBV3	25470	37401	1.33%	0.178%
28	14.057	1997	2001	2004	rBV	873642	757281	26.92%	3.614%
29	14.080	2004	2005	2011	rVB	42218	34108	1.21%	0.163%
30	14.151	2014	2017	2021	rVB2	29088	33361	1.19%	0.159%
31	14.804	2125	2128	2136	rBV	83209	104053	3.70%	0.497%
32	15.104	2176	2179	2183	rBV	21123	30165	1.07%	0.144%
33	15.480	2240	2243	2249	rVB2	24482	33990	1.21%	0.162%
34	15.545	2249	2254	2262	rVB2	472145	593066	21.08%	2.830%

Sum of corrected areas: 20955697

Instrument :

BNA_F

ClientSampleId :

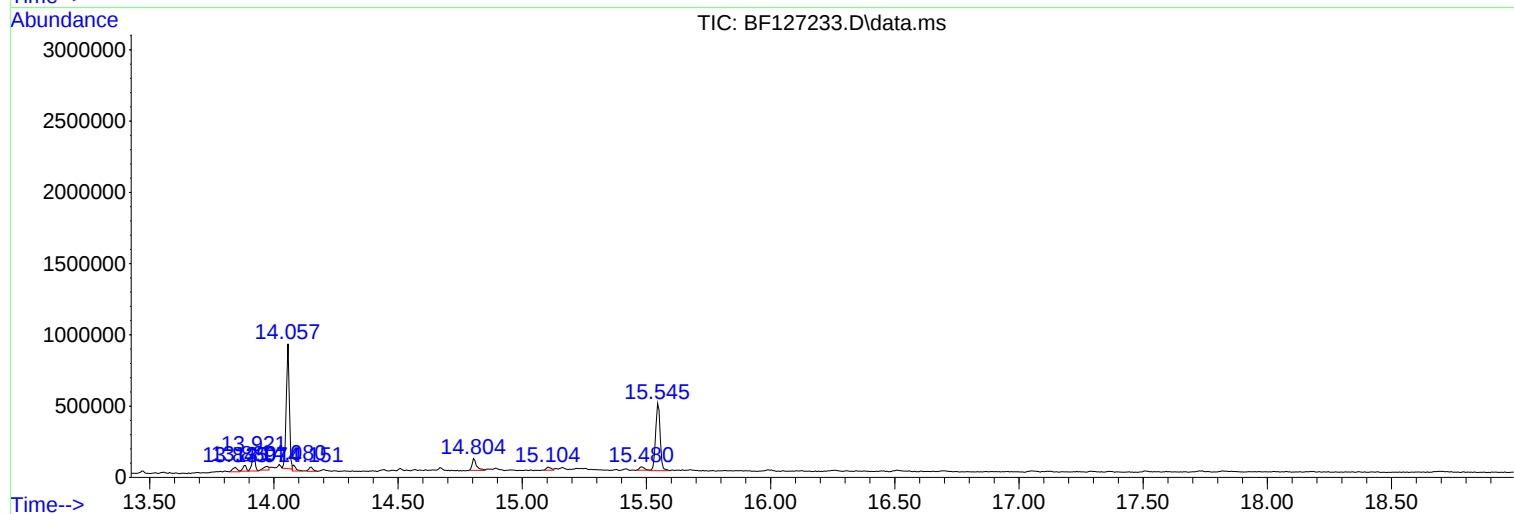
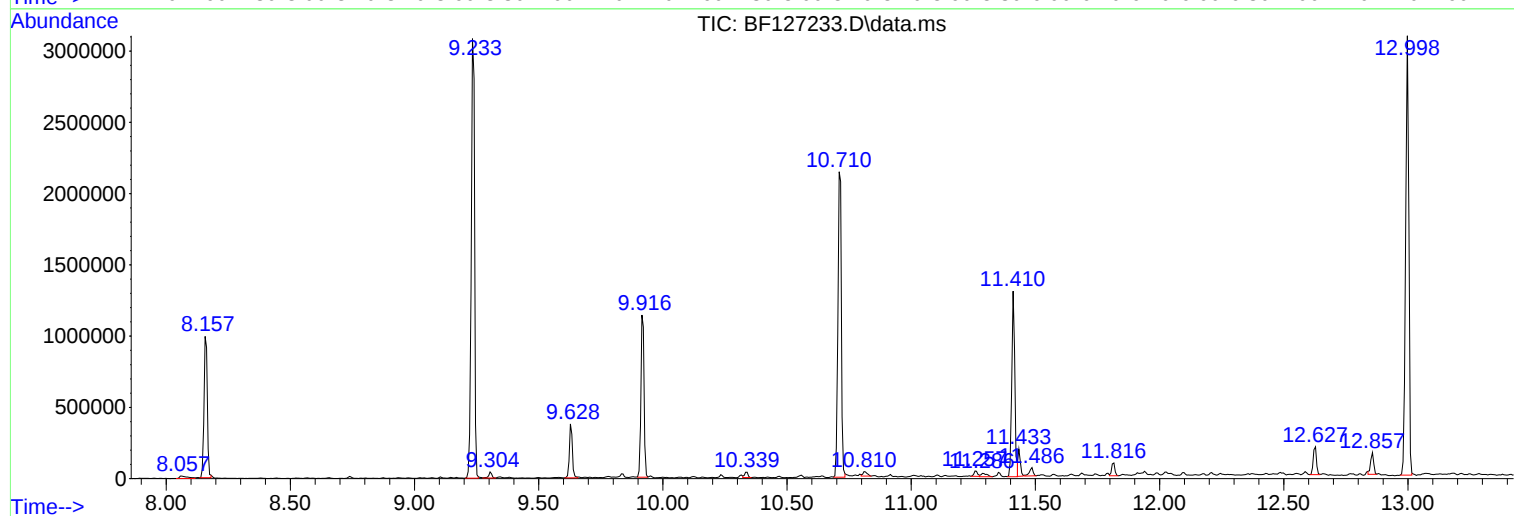
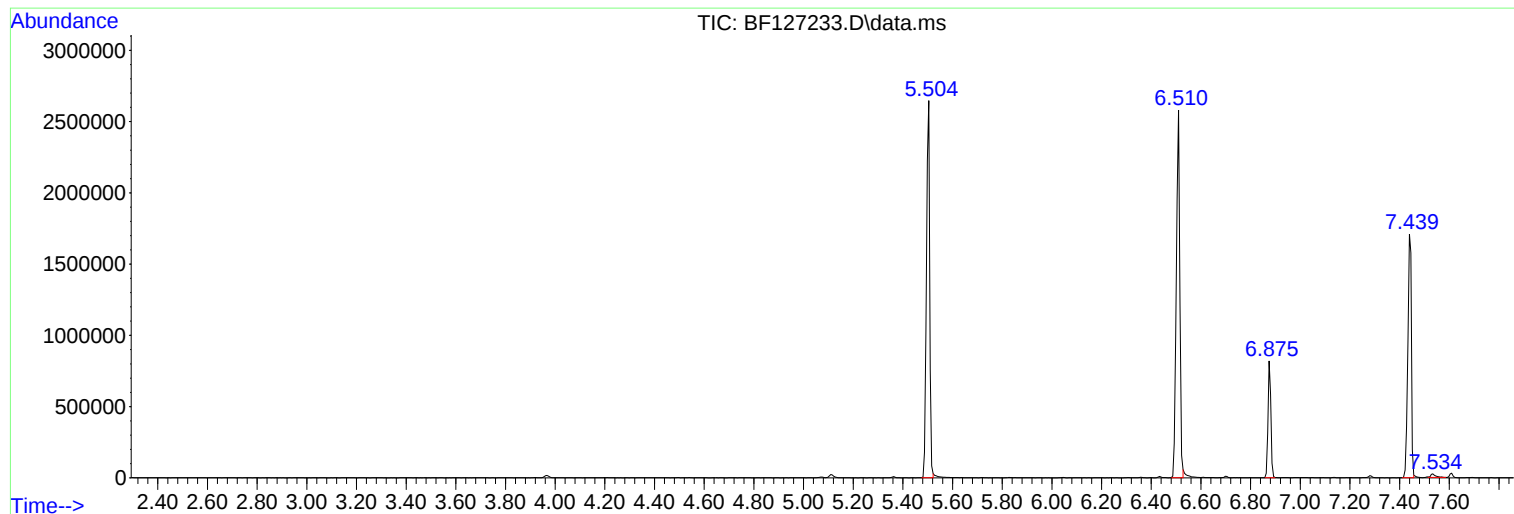
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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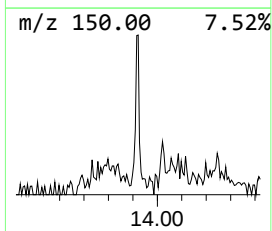
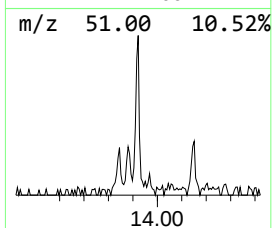
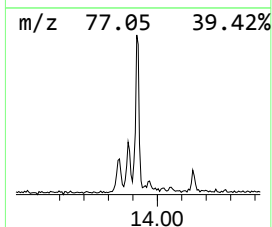
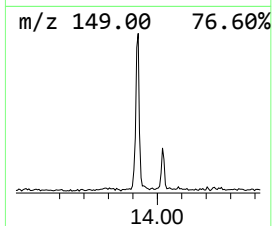
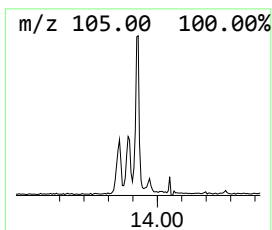
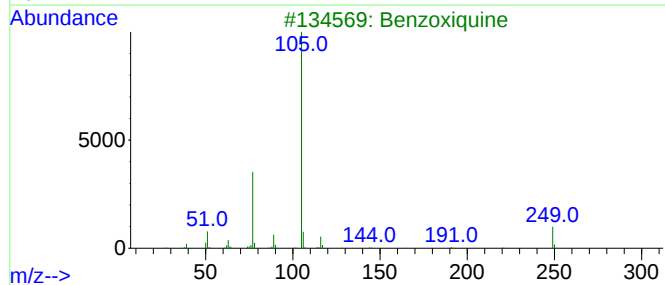
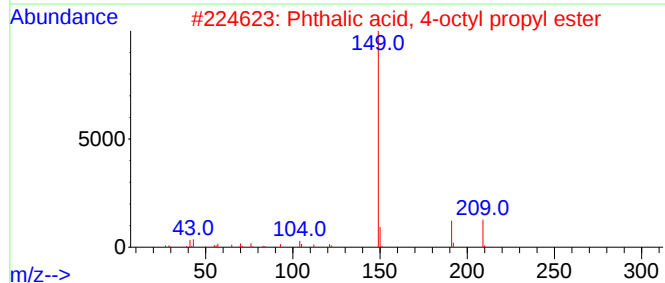
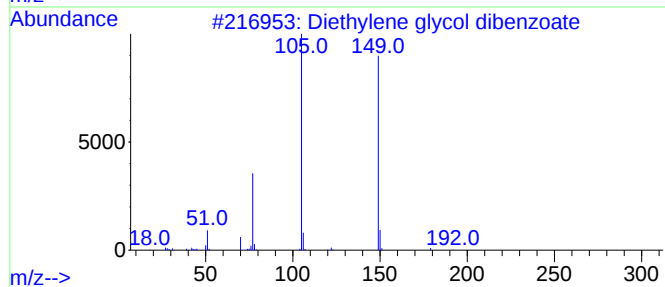
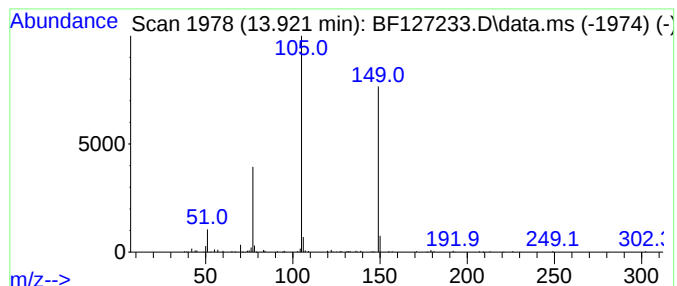
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Diethylene glycol dibenzoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.921	2.87 ng	108512	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2			Phthalic acid, 4-octyl propyl ester	320	C19H28O4	1000314-83-7	43
3			Benzoxiquine	249	C16H11NO2	000086-75-9	38
4			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	35
5			1,3-propanedione, 2-(acetyloxy)-...	282	C17H14O4	1000401-92-5	35



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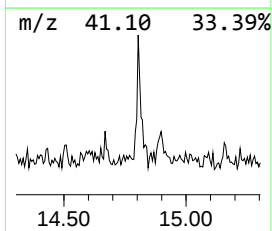
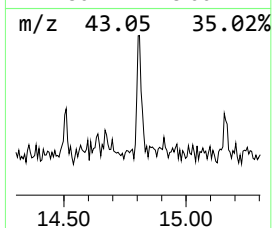
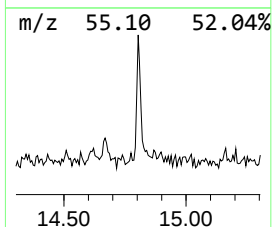
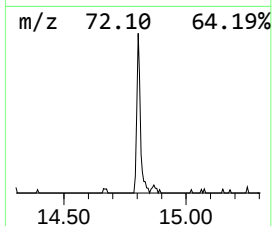
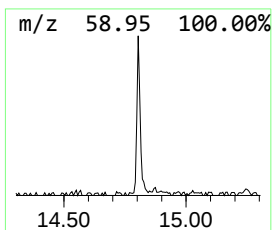
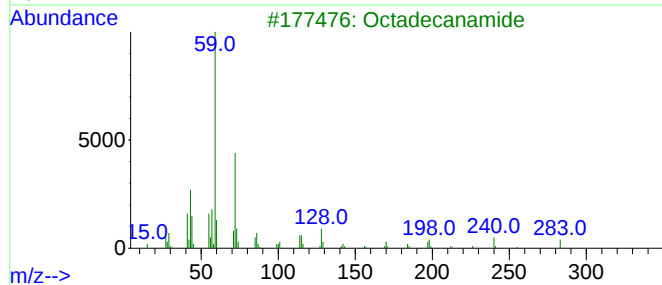
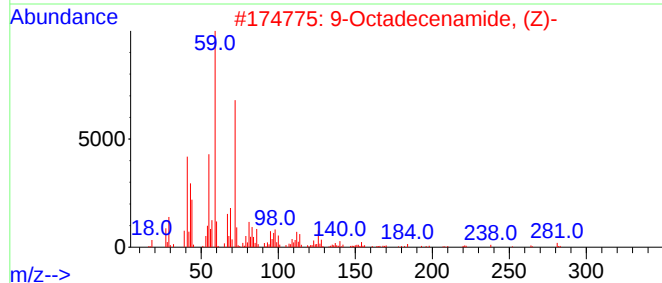
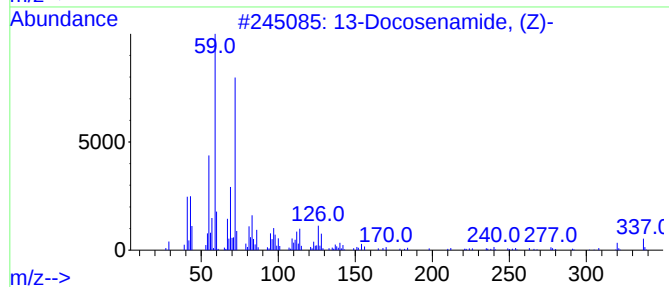
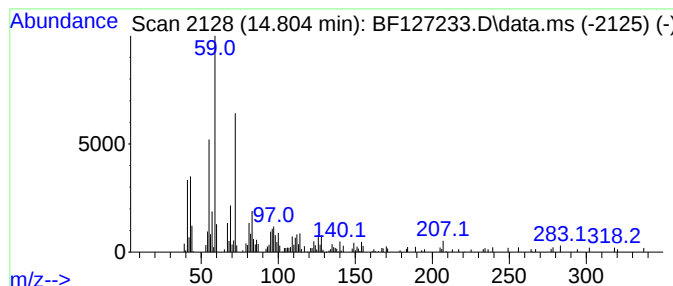
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Peak Number 2 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.804	3.51 ng	104053	Perylene-d12	15.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	87
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	86
3	Octadecanamide	283	C18H37NO	000124-26-5	83
4	Palmitoleamide	253	C16H31NO	106010-22-4	72
5	Dodecanamide	199	C12H25NO	001120-16-7	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Diethylene glyc...	13.921	2.9	ng	108512	5	14.057	757281	20.0
13-Docosenamide...	14.804	3.5	ng	104053	6	15.545	593066	20.0