

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121420\
 Data File : BF122711.D
 Acq On : 14 Dec 2020 17:56
 Operator : JU/CG
 Sample : L5046-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 11920

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-BF120520.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF122711.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.204	15	20	26	rVB3	38858	71381	1.84%	0.332%
2	5.340	549	553	565	rBV2	27657	70192	1.81%	0.326%
3	5.451	568	572	588	rVV	1064503	1094780	28.23%	5.090%
4	6.463	740	744	763	rBV2	700065	984686	25.39%	4.578%
5	6.834	804	807	816	rBV	1024035	816862	21.07%	3.798%
6	7.398	898	903	906	rBV	1924209	1778624	45.87%	8.269%
7	7.598	934	937	943	rVB	61651	51297	1.32%	0.238%
8	8.122	1022	1026	1029	rBV	1172337	1053014	27.15%	4.895%
9	8.151	1029	1031	1042	rVB2	57511	80393	2.07%	0.374%
10	9.086	1184	1190	1193	rBV2	30217	42953	1.11%	0.200%
11	9.198	1203	1209	1212	rBV	4206391	3877794	100.00%	18.028%
12	9.586	1271	1275	1279	rBV	235911	218034	5.62%	1.014%
13	9.875	1320	1324	1328	rBV	1525673	1296432	33.43%	6.027%
14	9.910	1328	1330	1335	rVB	50504	42621	1.10%	0.198%
15	10.663	1451	1458	1468	rBV	2074224	2077902	53.58%	9.660%
16	11.298	1564	1566	1571	rBV	43963	46057	1.19%	0.214%
17	11.363	1573	1577	1584	rVV	1283558	1272916	32.83%	5.918%
18	11.886	1663	1666	1667	rBV	69193	56579	1.46%	0.263%
19	12.451	1760	1762	1766	rVB2	51340	45554	1.17%	0.212%
20	12.674	1797	1800	1806	rBV5	62719	98330	2.54%	0.457%
21	12.945	1842	1846	1850	rBV	3651516	3782702	97.55%	17.585%
22	13.874	2001	2004	2010	rBV3	136707	213207	5.50%	0.991%
23	13.980	2019	2022	2023	rBV	321904	284723	7.34%	1.324%
24	13.998	2023	2025	2031	rVB	993862	977572	25.21%	4.545%
25	14.315	2077	2079	2083	rVB	141637	142127	3.67%	0.661%
26	15.462	2270	2274	2281	rBV2	754994	1033700	26.66%	4.806%

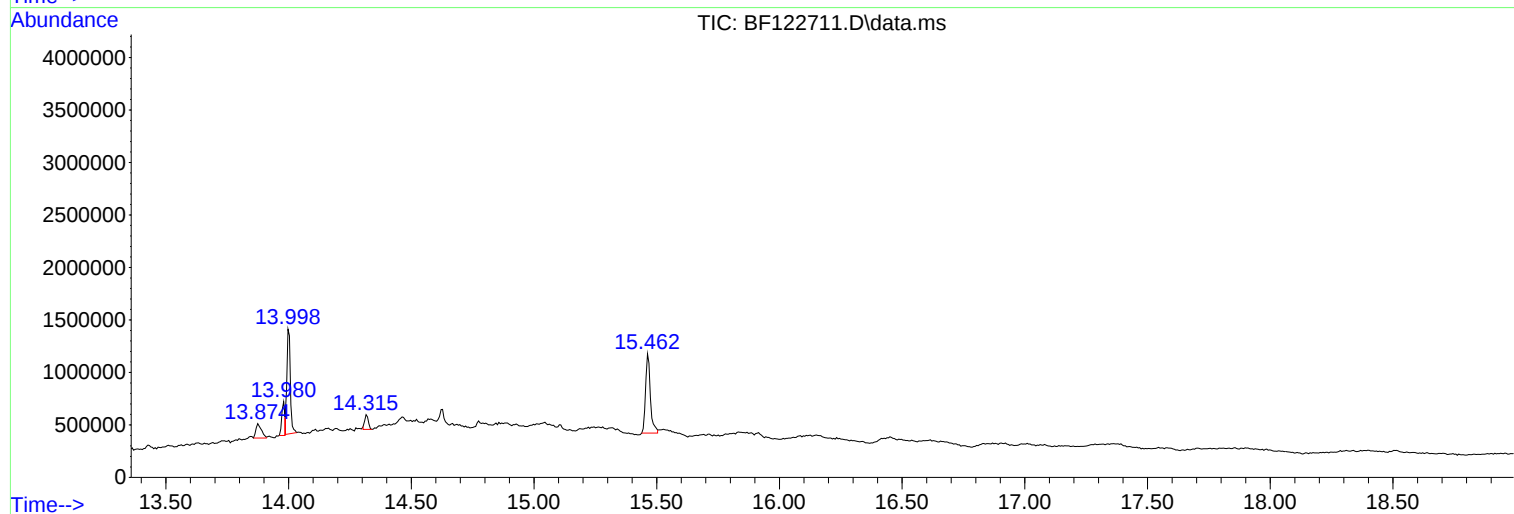
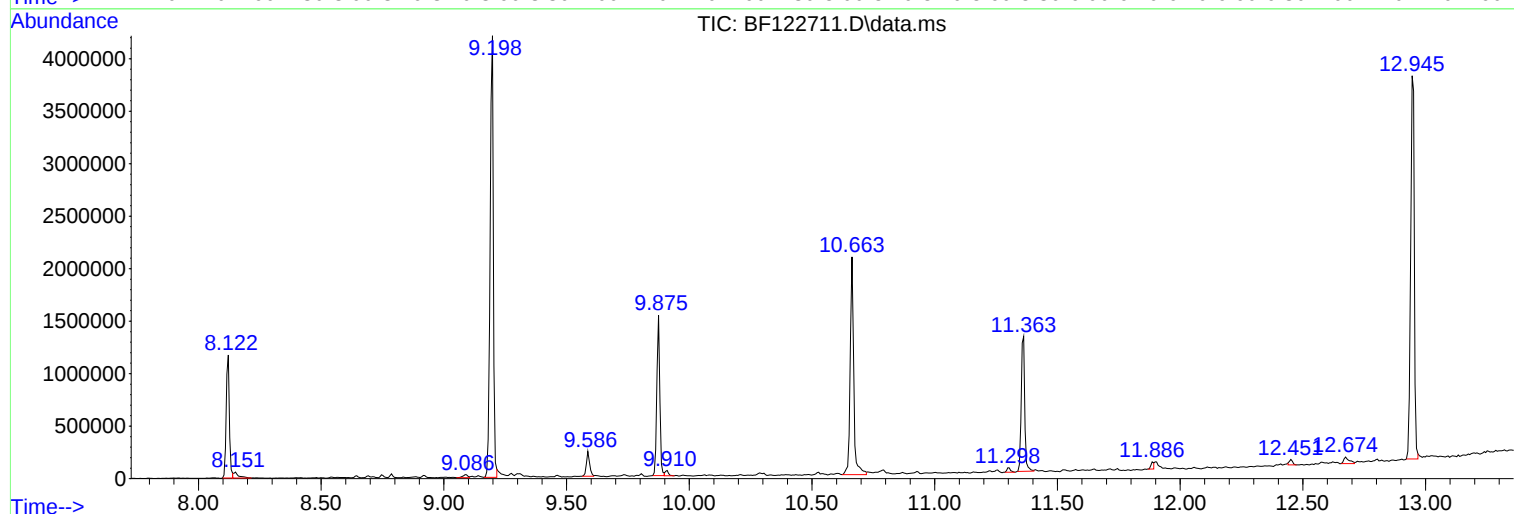
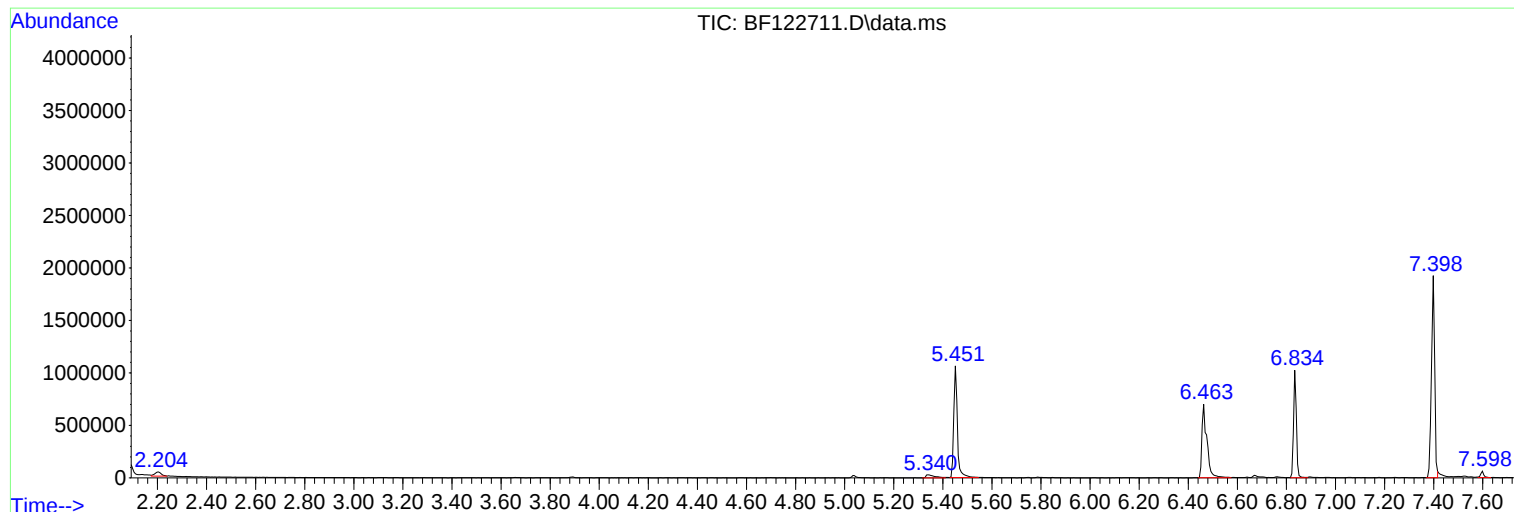
Sum of corrected areas: 21510432

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.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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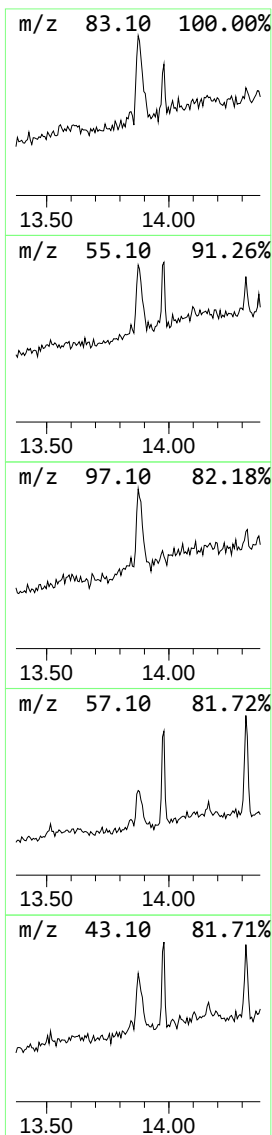
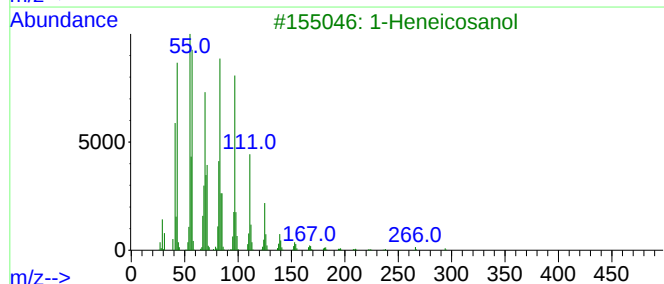
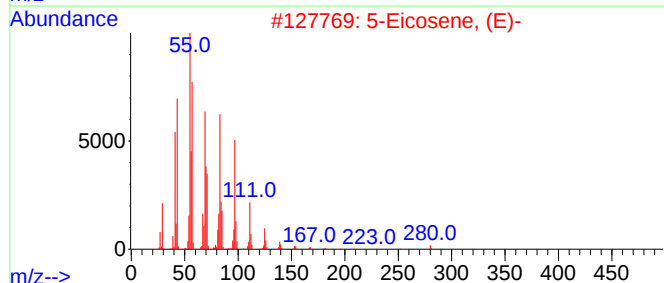
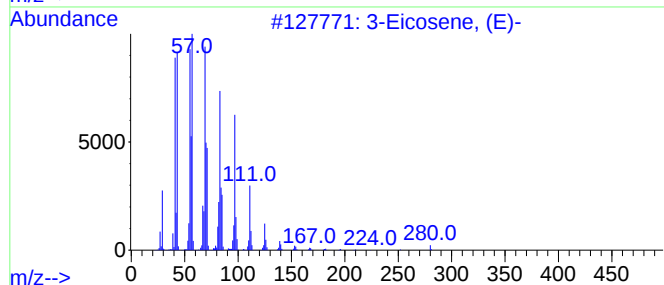
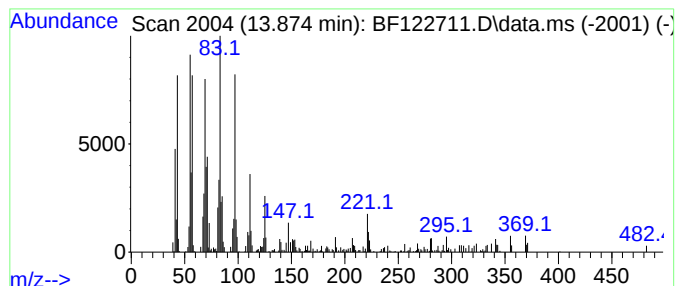
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Peak Number 1 3-Eicosene, (E)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	4.36 ng	213207	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Eicosene, (E)-	280	C20H40	074685-33-9	99
2			5-Eicosene, (E)-	280	C20H40	074685-30-6	98
3			1-Heneicosanol	312	C21H44O	015594-90-8	94
4			Behenic alcohol	326	C22H46O	000661-19-8	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	90



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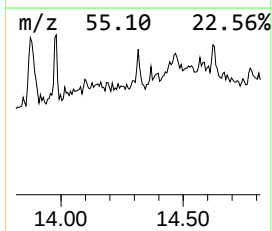
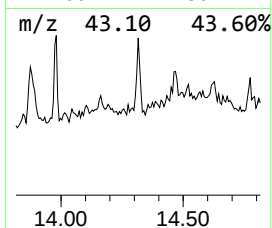
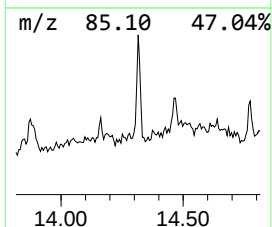
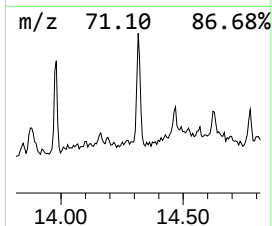
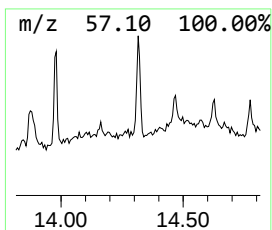
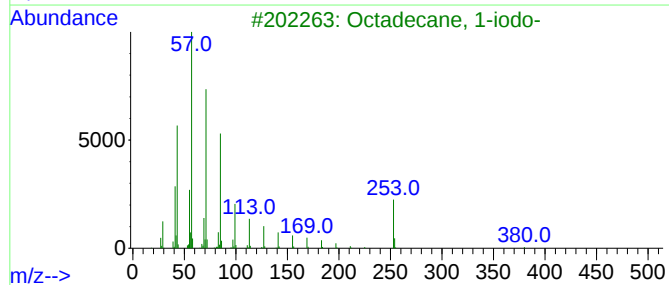
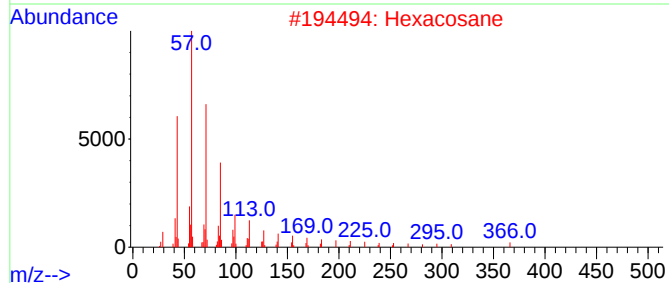
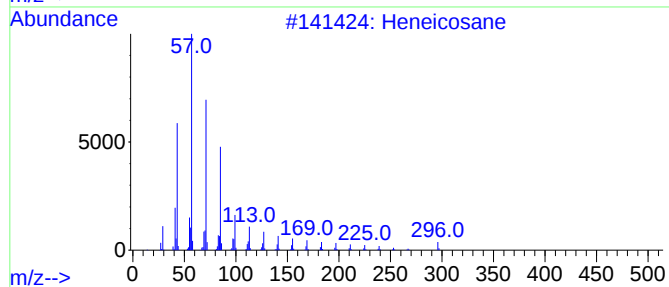
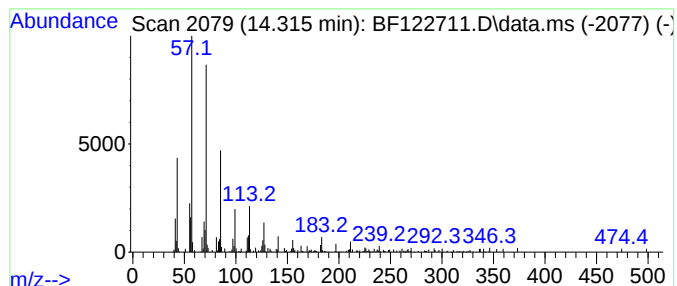
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Peak Number 2 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.315	2.91 ng	142127	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	93
2			Hexacosane	366	C26H54	000630-01-3	86
3			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
4			2-Bromotetradecane	276	C14H29Br	074036-95-6	86
5			Docosane, 11-decyl-	451	C32H66	055401-55-3	86



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Eicosene, (E)-	13.874	4.4	ng	213207	5	13.998	977572	20.0
Heneicosane	14.315	2.9	ng	142127	5	13.998	977572	20.0