

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113513.D
 Acq On : 2 Apr 2019 18:47
 Operator : JU/SJ
 Sample : K2179-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-5S

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF032519.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.304	543	547	571	rBV	2345328	2605170	88.04%	9.570%
2	6.339	719	723	740	rBV	2664593	2823350	95.42%	10.371%
3	6.463	740	744	755	rVB	3105647	2720520	91.94%	9.994%
4	6.692	779	783	793	rBV	914829	835529	28.24%	3.069%
5	6.845	805	809	825	rVB	2345938	2069243	69.93%	7.601%
6	7.257	874	879	892	rBV	1709573	1712687	57.88%	6.291%
7	7.975	997	1001	1016	rBV	1097331	1074712	36.32%	3.948%
8	8.457	1080	1083	1095	rBV	45371	99912	3.38%	0.367%
9	9.045	1179	1183	1186	rBV	3644640	2958980	100.00%	10.870%
10	9.439	1247	1250	1255	rBV	244686	202138	6.83%	0.743%
11	9.716	1294	1297	1307	rVB	1244290	1186422	40.10%	4.358%
12	10.504	1427	1431	1445	rBV	1914666	1936826	65.46%	7.115%
13	11.198	1545	1549	1556	rBV	1350782	1157102	39.10%	4.251%
14	11.733	1637	1640	1654	rBV	176116	236159	7.98%	0.868%
15	12.515	1770	1773	1784	rBV2	46822	65950	2.23%	0.242%
16	12.780	1813	1818	1821	rBV	2763070	2469316	83.45%	9.071%
17	13.686	1969	1972	1974	rBV2	50310	55834	1.89%	0.205%
18	13.827	1992	1996	2004	rVB	1086409	994002	33.59%	3.651%
19	13.998	2022	2025	2029	rVB	60282	50697	1.71%	0.186%
20	14.304	2065	2077	2088	rVB6	114518	310915	10.51%	1.142%
21	14.598	2123	2127	2130	rBV	95592	103649	3.50%	0.381%
22	14.662	2135	2138	2143	rVB4	32610	41121	1.39%	0.151%
23	14.904	2176	2179	2184	rVB	95786	101229	3.42%	0.372%
24	15.233	2230	2235	2246	rBV	910860	1202122	40.63%	4.416%
25	15.645	2301	2305	2310	rBV	74382	108606	3.67%	0.399%
26	16.103	2379	2383	2395	rVB	53930	100156	3.38%	0.368%

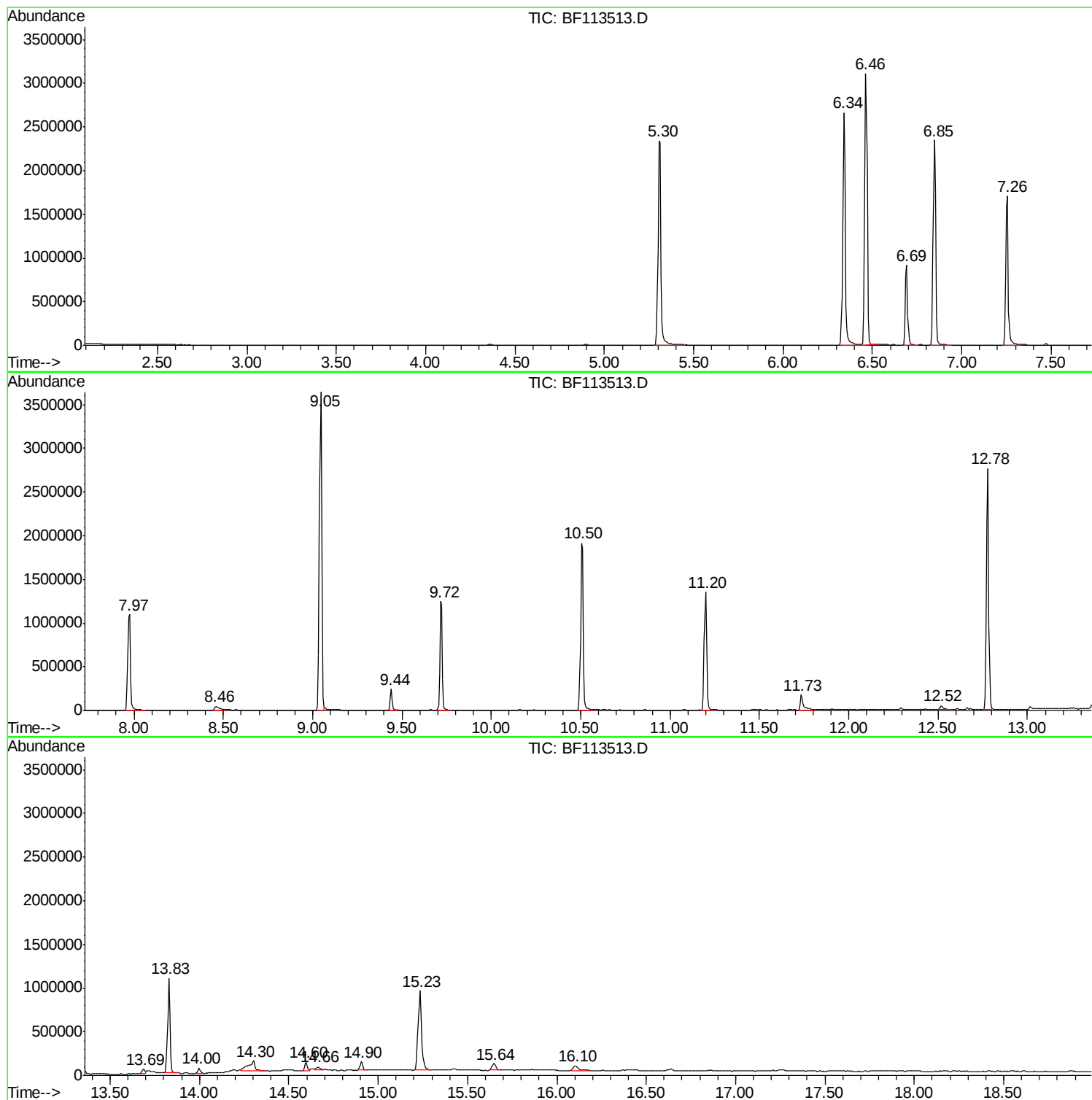
Sum of corrected areas: 27222347

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.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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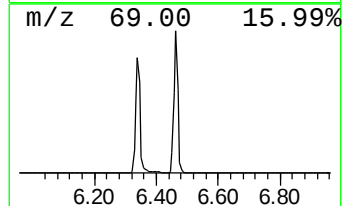
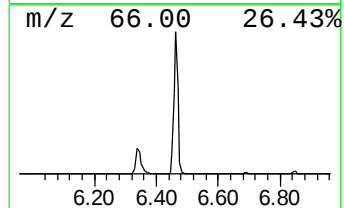
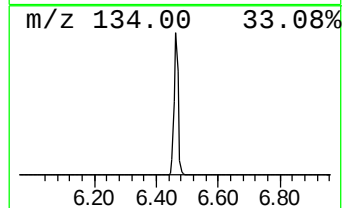
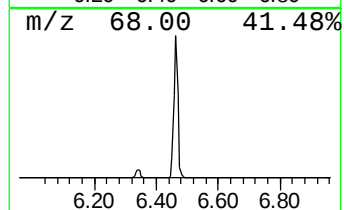
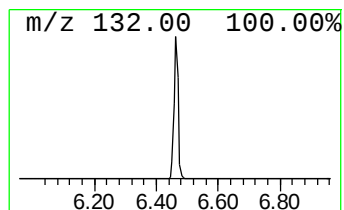
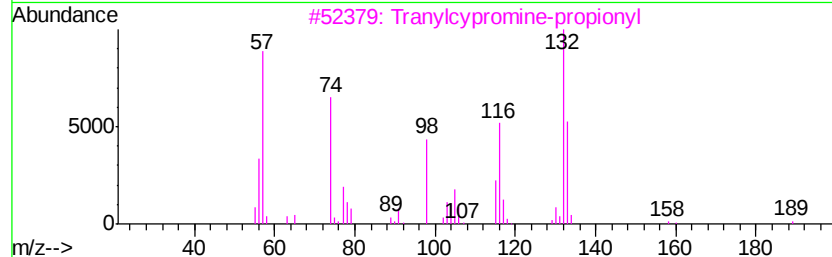
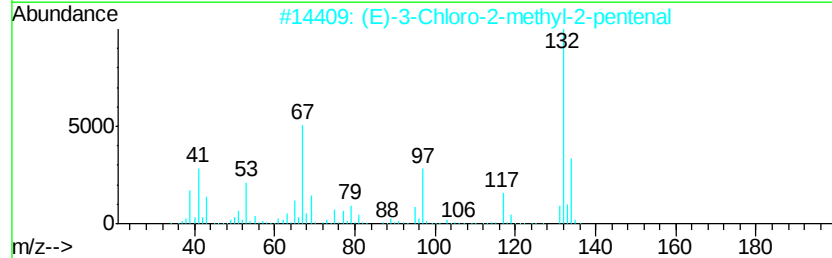
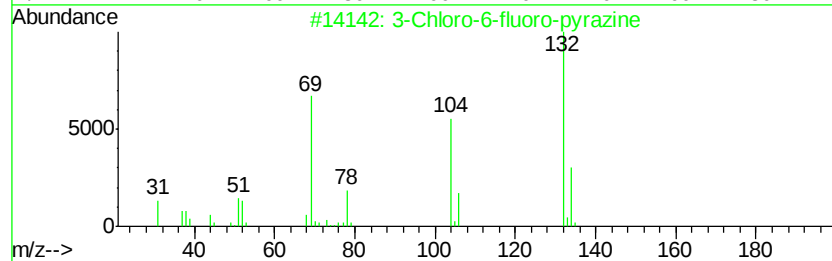
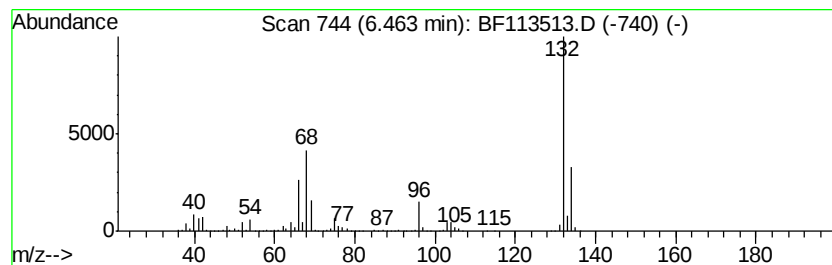
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	65.12 ng	2720520	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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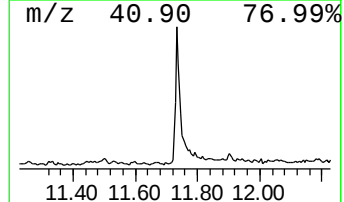
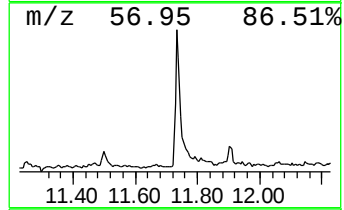
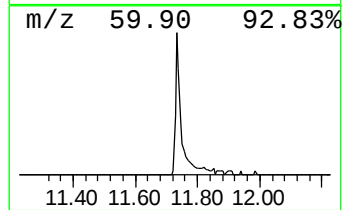
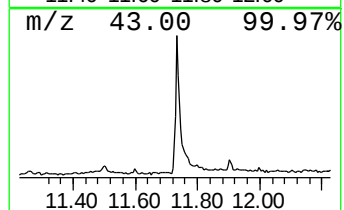
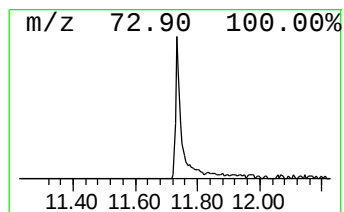
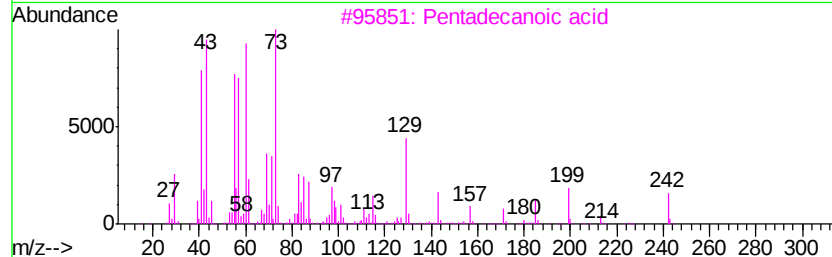
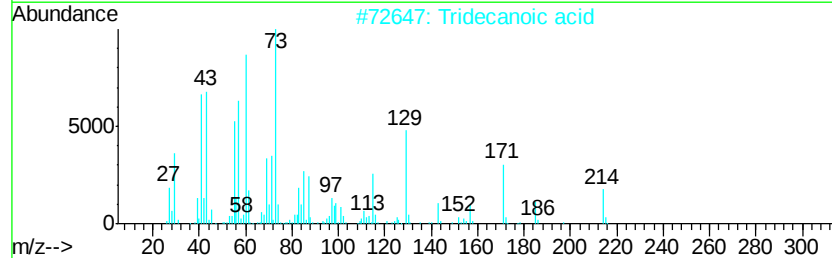
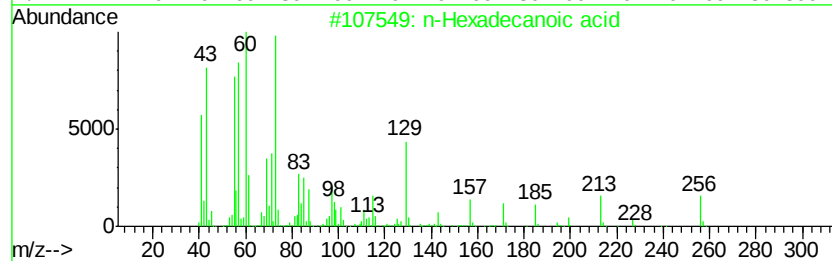
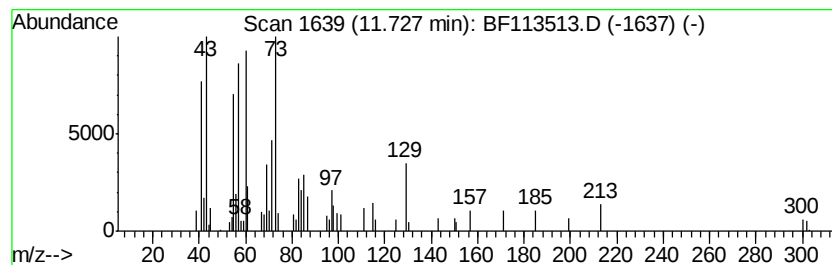
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	4.08 ng	236159	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Octadecanoic acid	284	C18H36O2	000057-11-4	90
5		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	30



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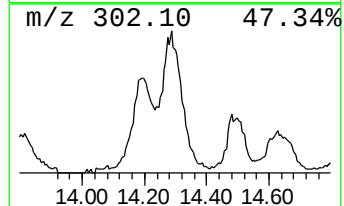
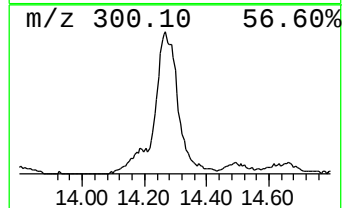
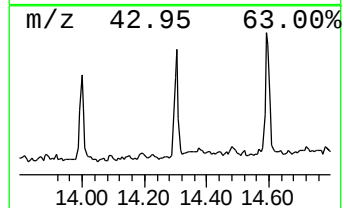
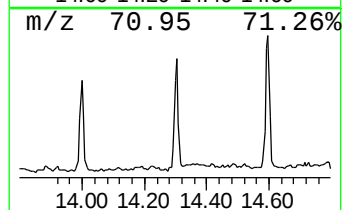
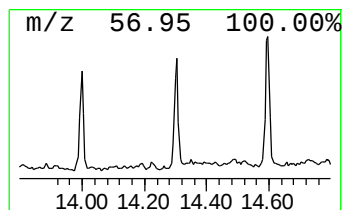
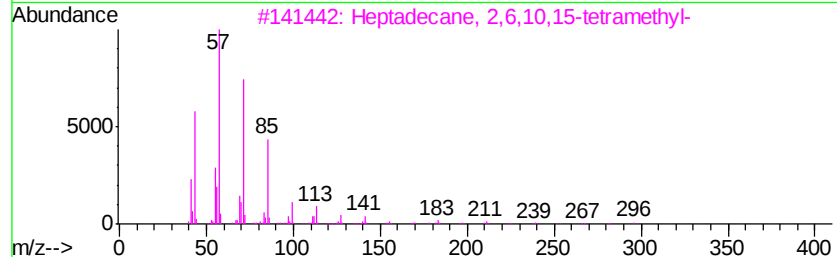
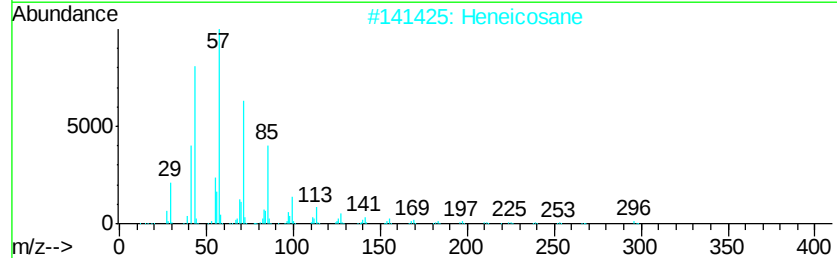
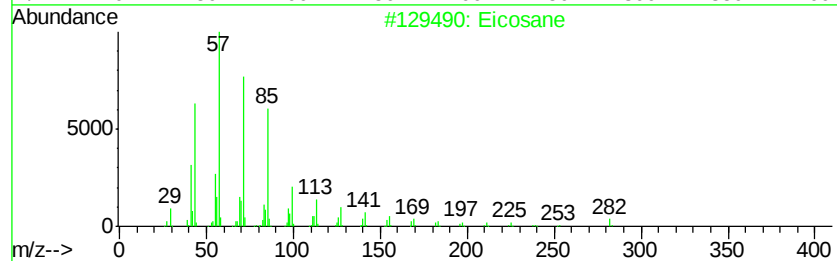
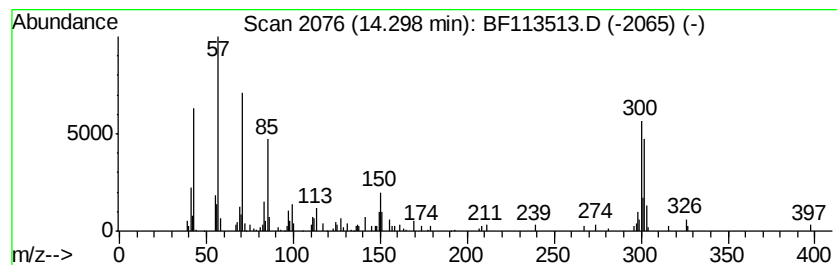
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Peak Number 4 Eicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	6.26 ng	310915	Chrysene-d12	13.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	89
2	Heneicosane		296	C21H44	000629-94-7	70
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	70
4	Triacontane		422	C30H62	000638-68-6	49
5	Tetracosane		338	C24H50	000646-31-1	49



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.46	6.46	65.1	ng	2720520	1	6.69	835529	20.0
n-Hexadecanoic acid	11.73	4.1	ng	236159	4	11.20	1157100	20.0
Eicosane	14.30	6.3	ng	310915	5	13.83	994002	20.0