

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100418\
 Data File : BF109601.D
 Acq On : 4 Oct 2018 21:36
 Operator : JU/SJ
 Sample : J5244-04 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-WATER

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-BF092218.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	5.316	546	549	562	rBV	16735	31675	4.28%	0.599%
2	6.357	721	726	735	rBV3	10284	27329	3.70%	0.517%
3	6.475	743	746	761	rBV	90660	133342	18.03%	2.522%
4	6.704	781	785	799	rBV	424159	494012	66.80%	9.342%
5	6.857	807	811	827	rVB	148445	149523	20.22%	2.828%
6	7.263	877	880	897	rBV	50718	105724	14.30%	1.999%
7	7.981	999	1002	1019	rBV	608929	678875	91.79%	12.838%
8	9.057	1182	1185	1202	rBV	225018	233888	31.63%	4.423%
9	9.451	1249	1252	1262	rBV	25519	25483	3.45%	0.482%
10	9.733	1296	1300	1318	rBV	765871	739560	100.00%	13.986%
11	10.151	1367	1371	1374	rBV	298865	243232	32.89%	4.600%
12	10.516	1430	1433	1445	rBV	66413	84976	11.49%	1.607%
13	11.210	1547	1551	1567	rBV	639883	621878	84.09%	11.760%
14	12.786	1816	1819	1828	rVB	179745	163692	22.13%	3.096%
15	13.698	1971	1974	1982	rBV5	7587	14110	1.91%	0.267%
16	13.833	1994	1997	2010	rVB	561046	558832	75.56%	10.568%
17	14.280	2069	2073	2076	rBV2	22215	26901	3.64%	0.509%
18	14.316	2076	2079	2082	rVB2	14803	13405	1.81%	0.253%
19	14.610	2125	2129	2134	rVB	35991	33978	4.59%	0.643%
20	14.921	2178	2182	2187	rBV	44172	49193	6.65%	0.930%
21	15.239	2231	2236	2240	rBV	481713	584258	79.00%	11.049%
22	15.268	2240	2241	2248	rVB	68788	69324	9.37%	1.311%
23	15.668	2304	2309	2316	rVB	60929	92107	12.45%	1.742%
24	16.127	2383	2387	2393	rVB2	46041	71151	9.62%	1.346%
25	16.668	2473	2479	2485	rBV3	20634	41533	5.62%	0.785%

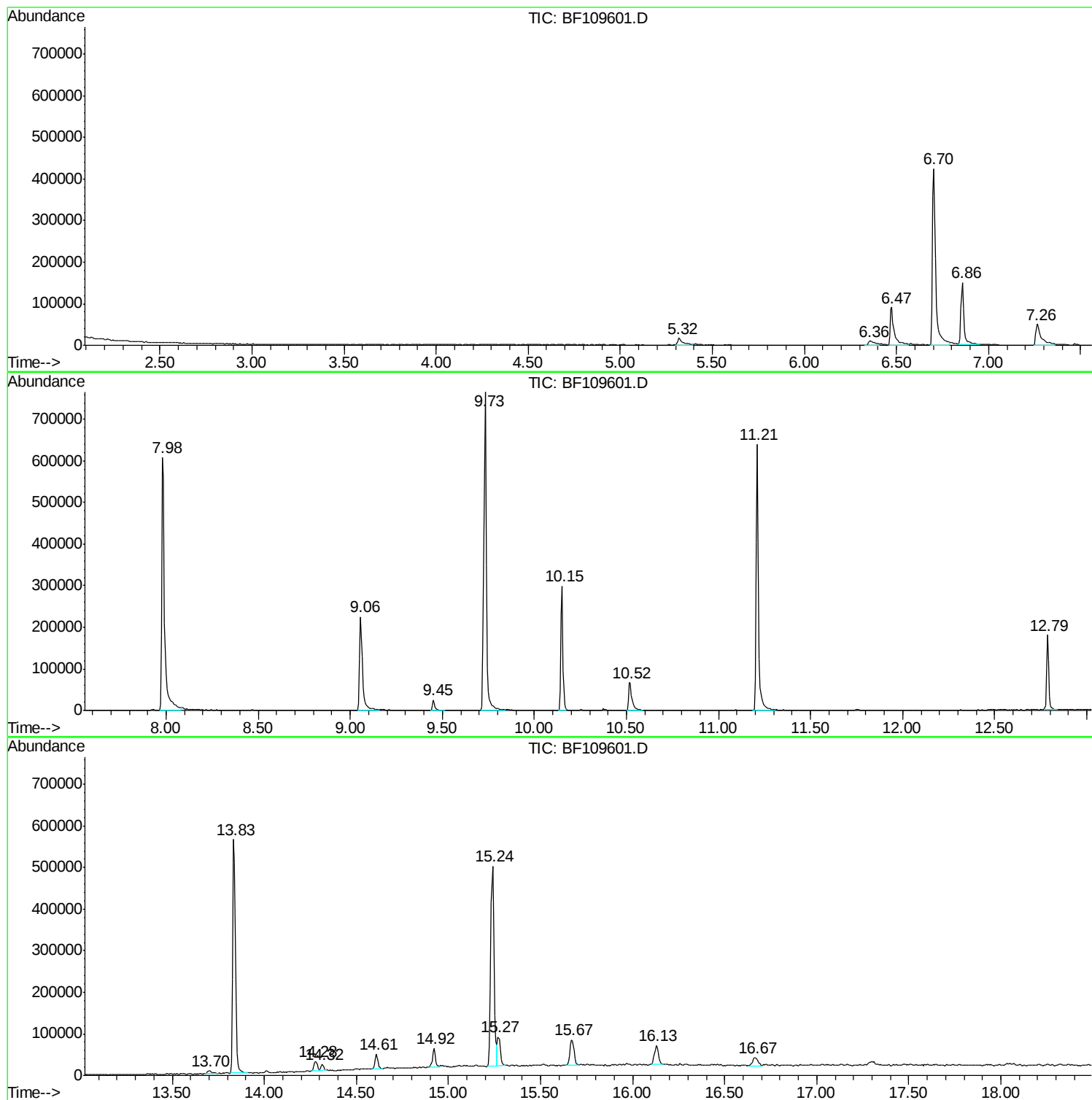
Sum of corrected areas: 5287981

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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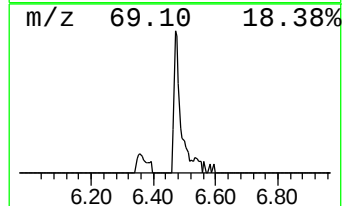
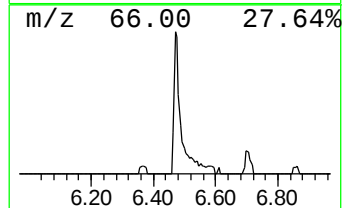
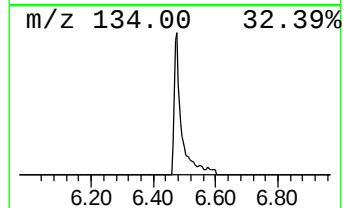
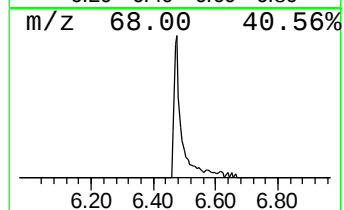
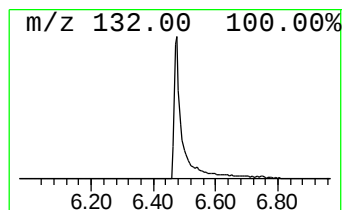
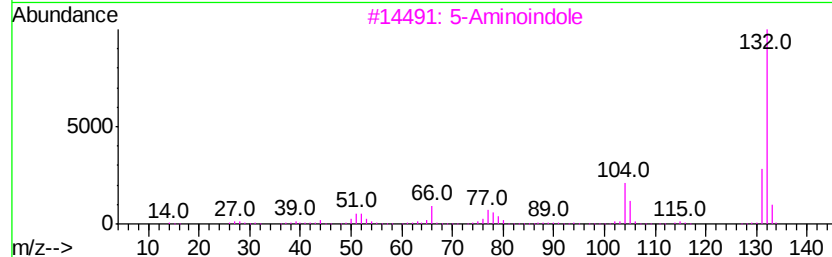
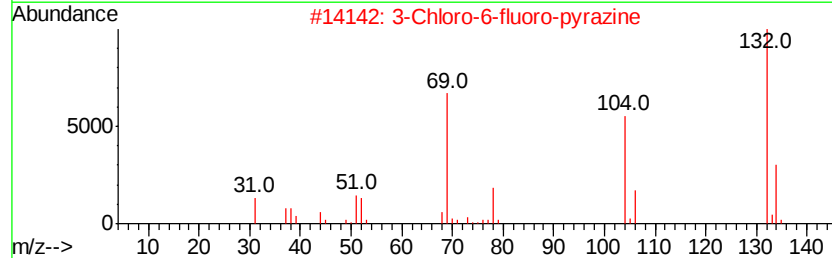
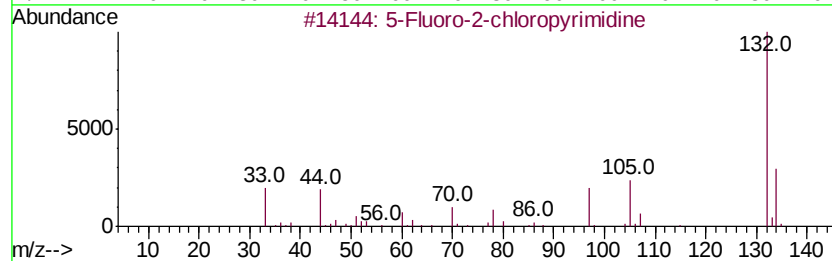
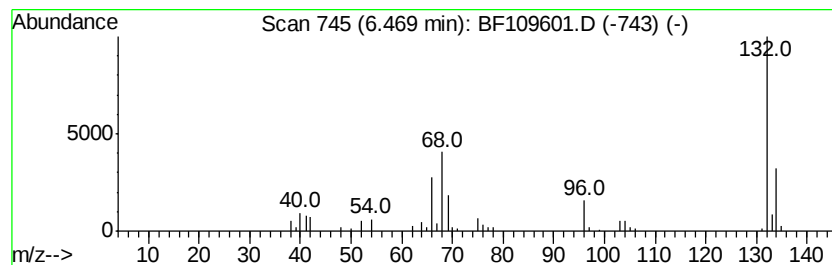
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	5.40 ng	133342	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3			5-Aminoindole	132	C8H8N2	005192-03-0	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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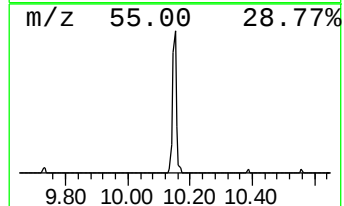
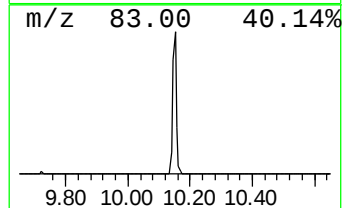
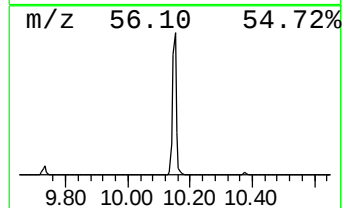
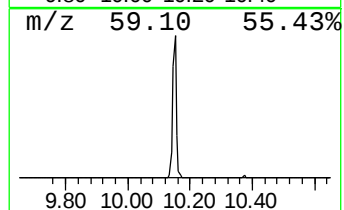
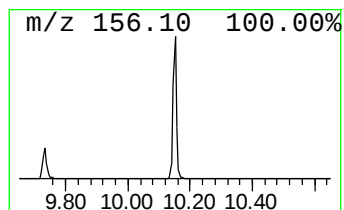
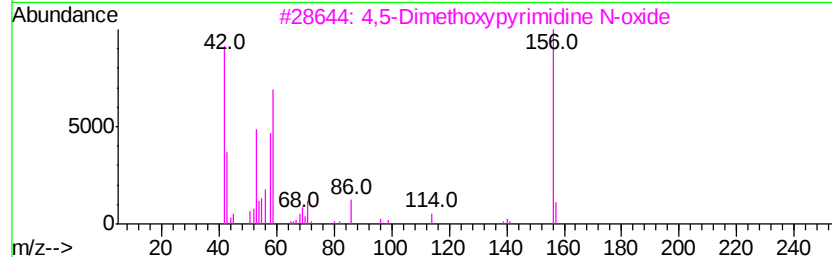
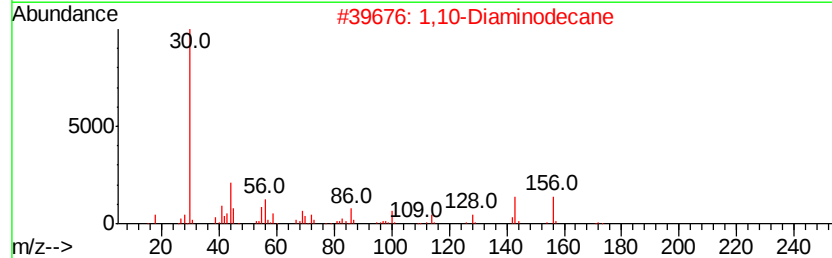
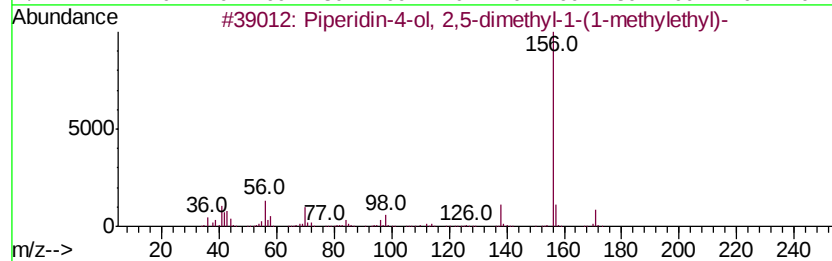
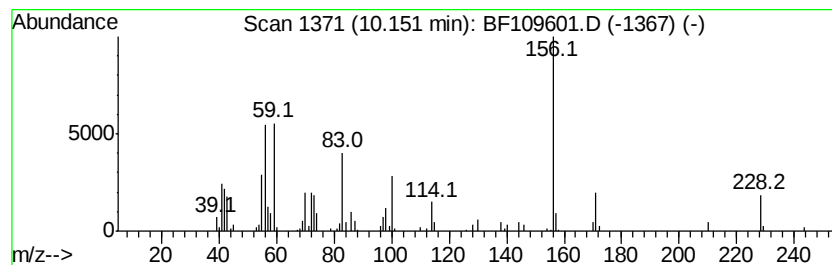
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Peak Number 3 unknown10.15 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.15	6.58 ng	243232	Acenaphthene-d10	9.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Piperidin-4-ol, 2,5-dimethyl-1-(...	171	C10H21NO	1000271-57-8	43
2		1,10-Diaminodecane	172	C10H24N2	000646-25-3	38
3		4,5-Dimethoxyxpyrimidine N-oxide	156	C6H8N2O3	114969-59-4	35
4		Oxalic acid, monoamide, N,N-dibu...	383	C23H45NO3	1000309-46-7	27
5		(S)-(-)-.alpha.-(1-Naphthyl)ethy...	171	C12H13N	010420-89-0	22



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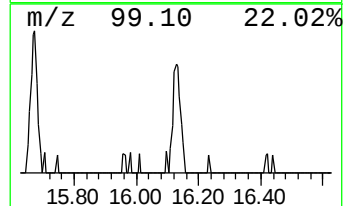
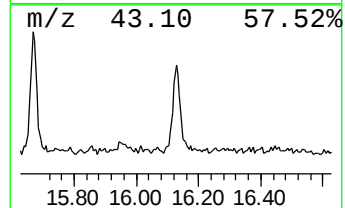
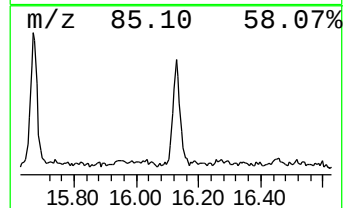
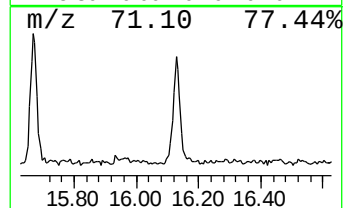
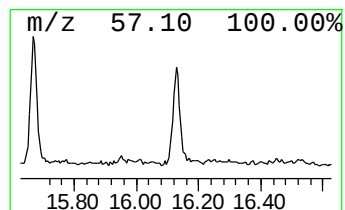
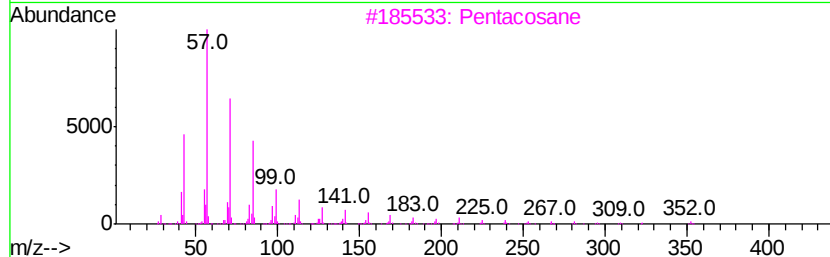
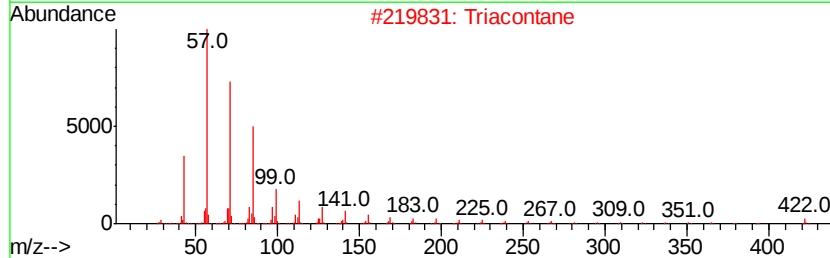
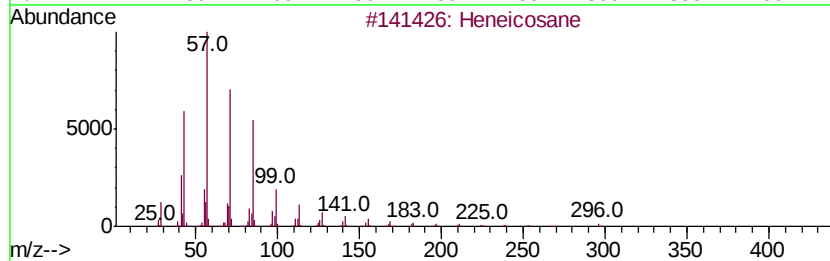
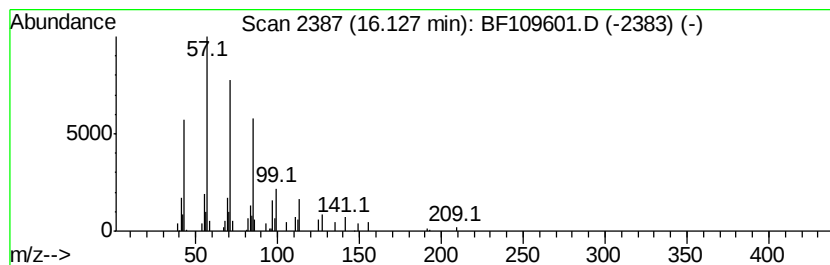
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Peak Number 5 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	2.44 ng	71151	Perylene-d12	15.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Triacontane		422	C30H62	000638-68-6	90
3	Pentacosane		352	C25H52	000629-99-2	87
4	Hexacosane		366	C26H54	000630-01-3	83
5	Octacosane		394	C28H58	000630-02-4	80



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.47	6.47	5.4	ng	133342	1	6.70	494012	20.0
unknown10.15	10.15	6.6	ng	243232	3	9.73	739560	20.0
Heneicosane	16.13	2.4	ng	71151	6	15.24	584258	20.0