

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
 Data File : BF081443.D
 Acq On : 5 Sep 2015 3:34
 Operator : UM/IZ
 Sample : G3511-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08(7-8)

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:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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	2.753	99	102	105	rBB	16049	28554	1.30%	0.159%
2	3.541	169	171	179	rBB	11998	26278	1.20%	0.147%
3	4.410	243	247	259	rVB	690707	1307480	59.48%	7.295%
4	4.913	288	291	293	rBV	1841468	2015406	91.68%	11.245%
5	5.336	326	328	334	rVB	25939	25314	1.15%	0.141%
6	6.033	386	389	391	rBV	1965132	2198269	100.00%	12.265%
7	6.124	395	397	400	rBV	1878082	1755330	79.85%	9.794%
8	6.353	415	417	420	rVB	366009	358327	16.30%	1.999%
9	6.513	428	431	433	rBV	1639705	1419409	64.57%	7.920%
10	6.936	465	468	470	rBV	1317927	1281745	58.31%	7.151%
11	7.645	528	530	533	rBV	553396	468889	21.33%	2.616%
12	8.730	622	625	627	rBV	2254996	2080712	94.65%	11.609%
13	9.130	658	660	663	rBV	327402	288149	13.11%	1.608%
14	9.382	680	682	685	rVB	474552	465845	21.19%	2.599%
15	10.171	748	751	753	rBV	1269684	1258757	57.26%	7.023%
16	10.319	762	764	767	rBV	32882	36168	1.65%	0.202%
17	10.765	801	803	804	rBV	29599	24875	1.13%	0.139%
18	10.845	808	810	813	rBV	414765	472280	21.48%	2.635%
19	11.428	859	861	864	rBV	36190	46439	2.11%	0.259%
20	12.434	946	949	952	rBV	1670067	1619818	73.69%	9.038%
21	13.359	1028	1030	1036	rBV	60634	128362	5.84%	0.716%
22	13.451	1036	1038	1041	rBV	334528	350610	15.95%	1.956%
23	14.765	1150	1153	1155	rBV	226980	265943	12.10%	1.484%

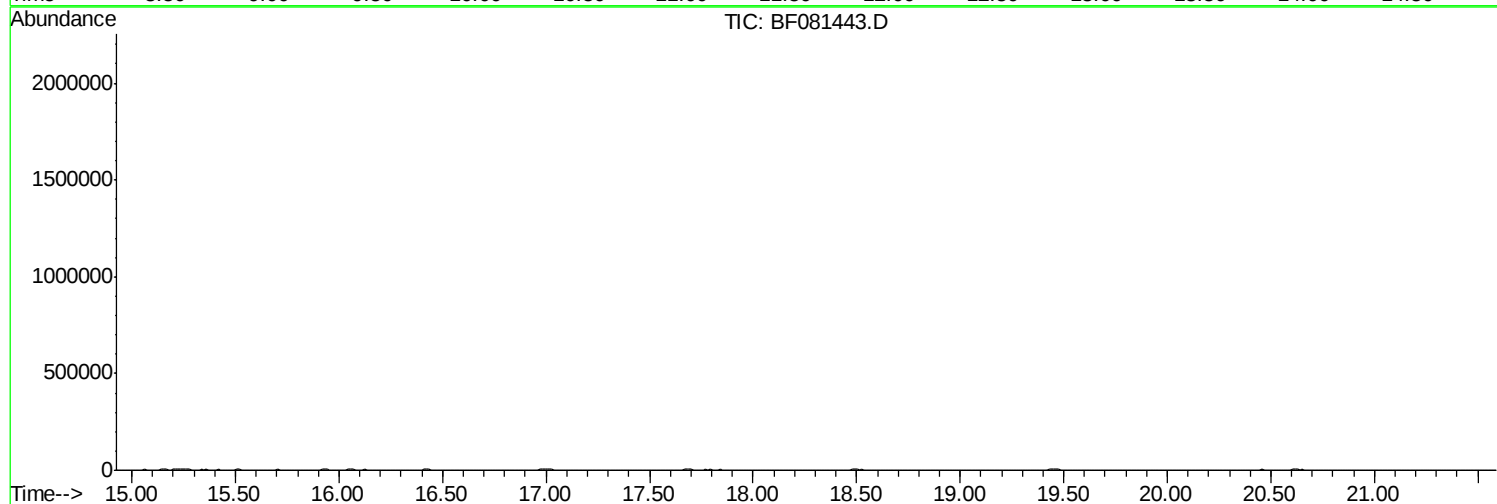
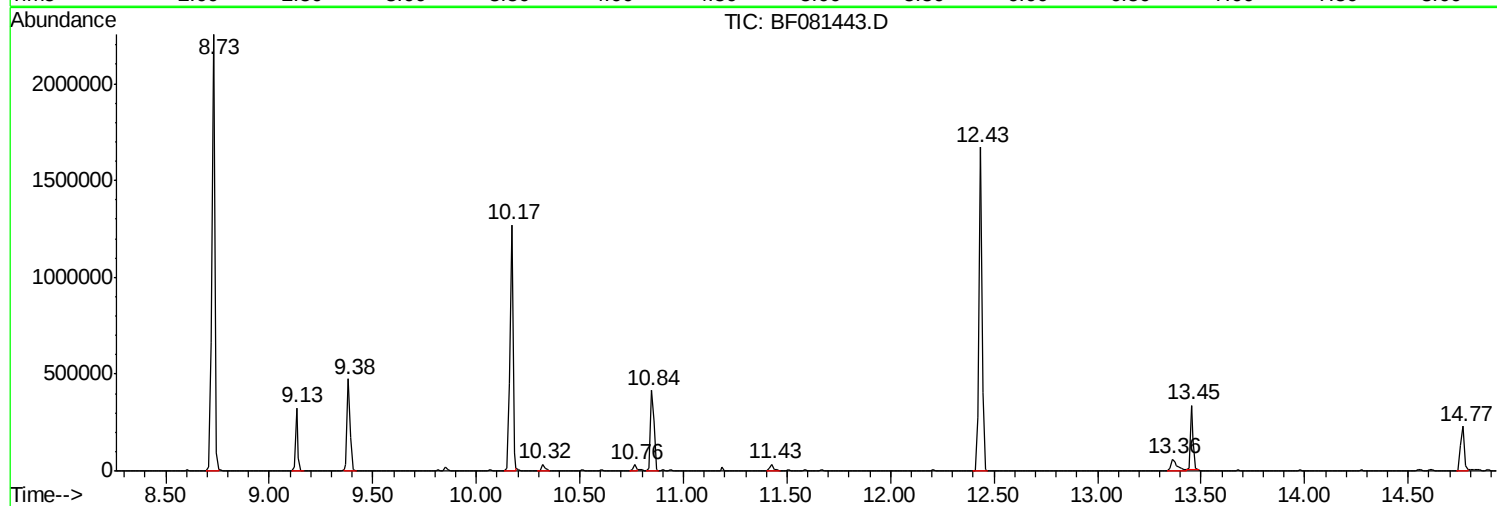
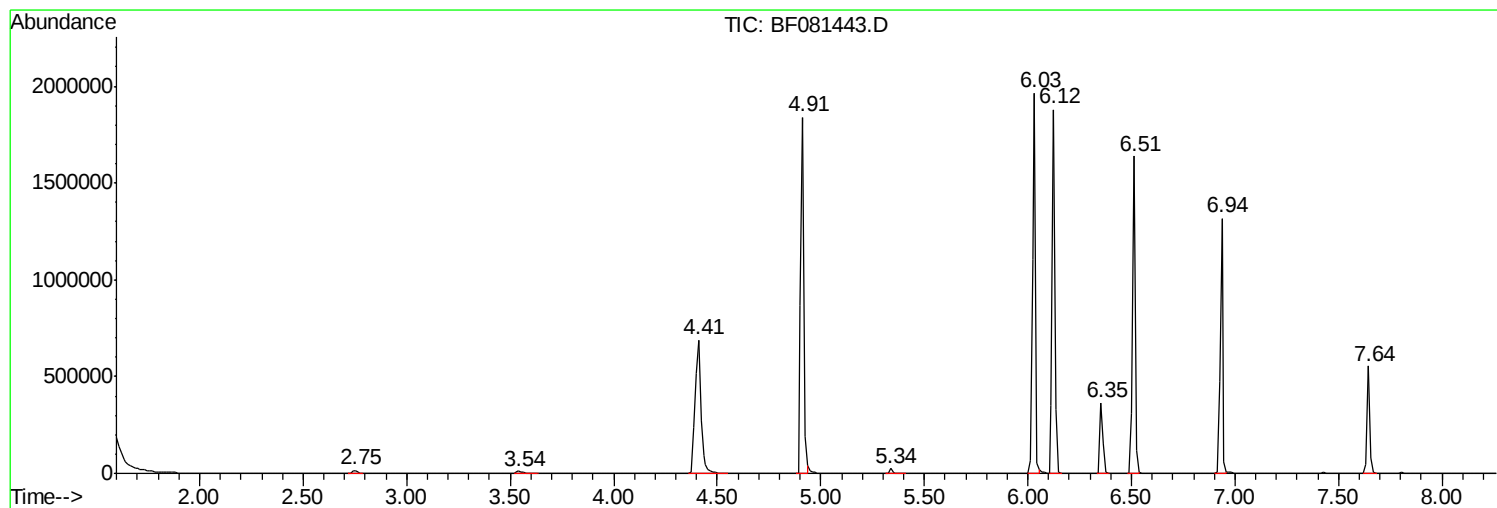
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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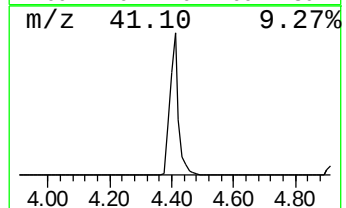
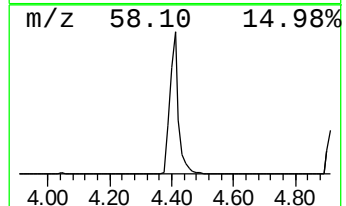
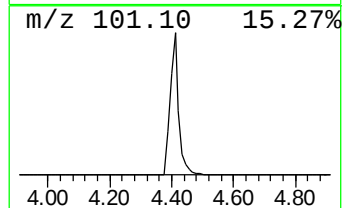
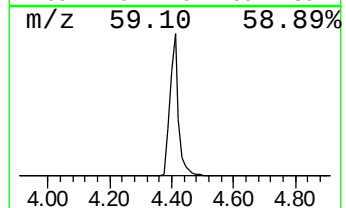
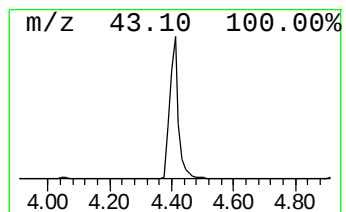
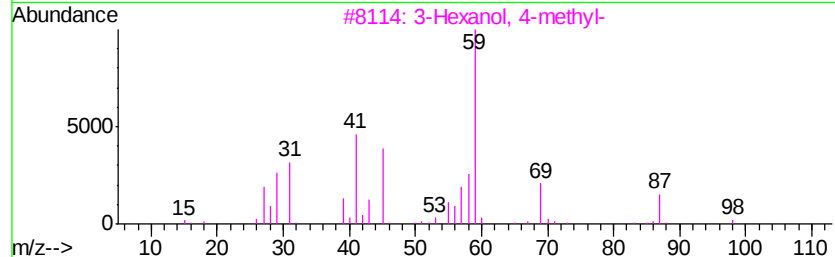
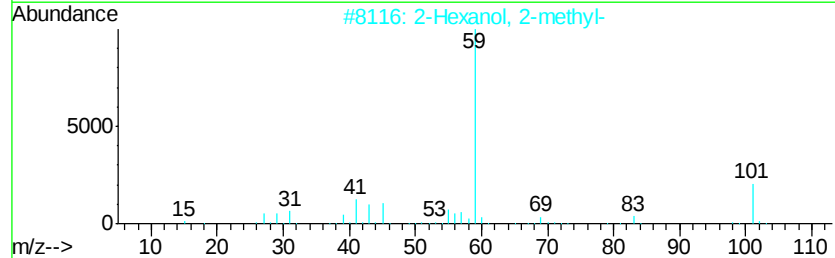
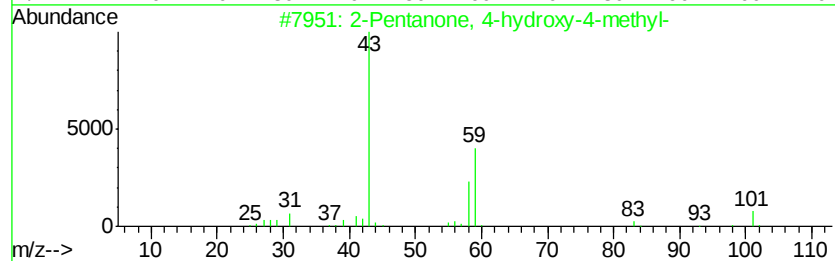
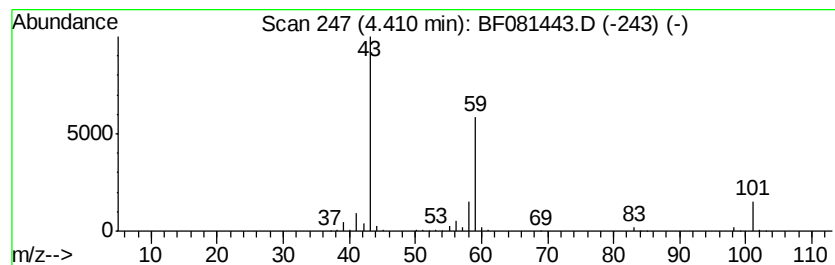
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.41	72.98 ng	1307480	1,4-Dichlorobenzene-d4	6.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17
5			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9



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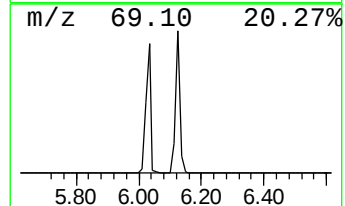
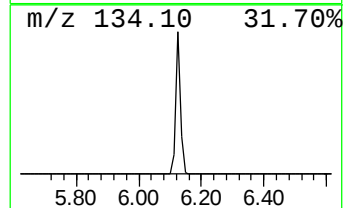
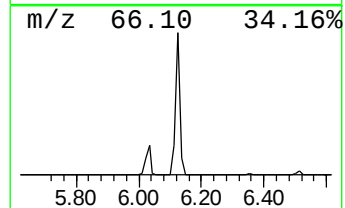
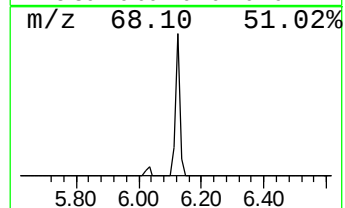
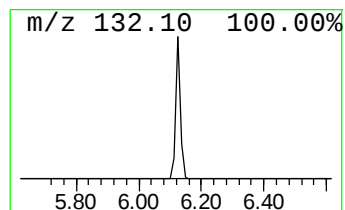
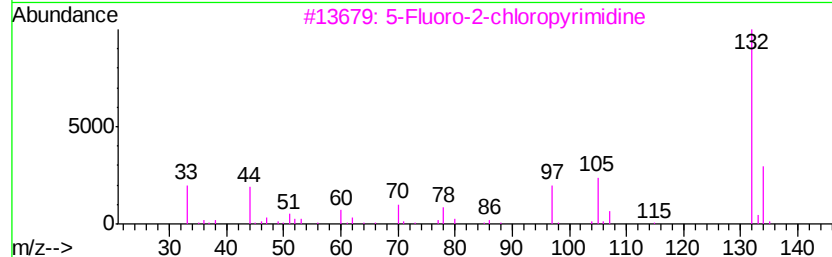
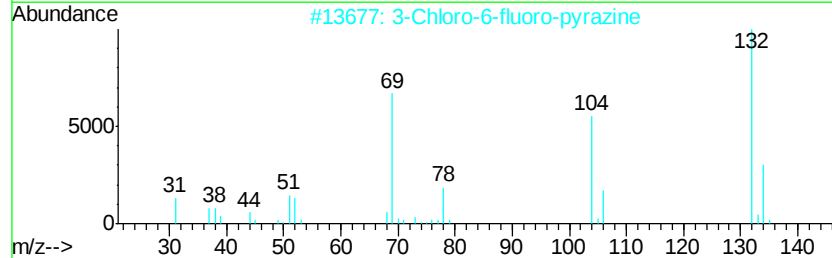
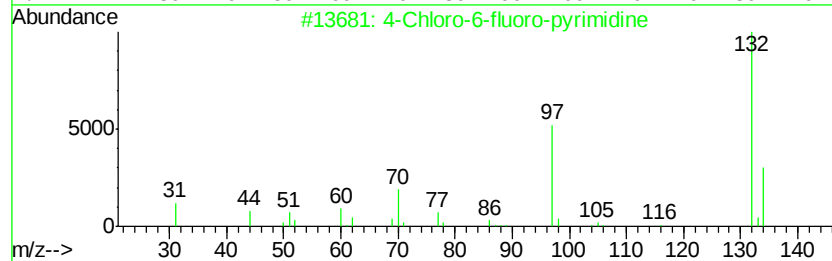
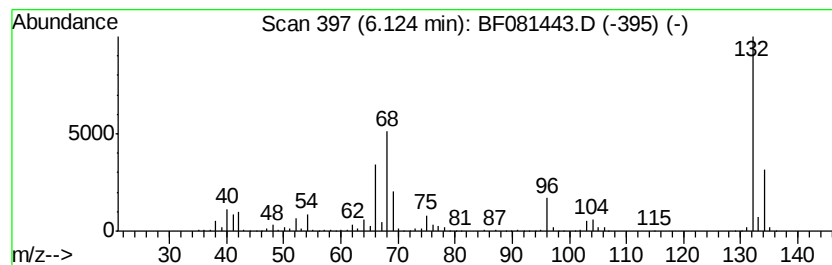
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Peak Number 2 unknown6.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	97.97 ng	1755330	1,4-Dichlorobenzene-d4	6.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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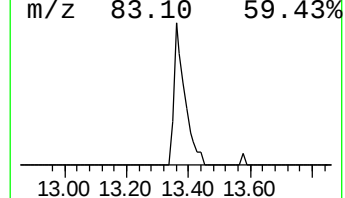
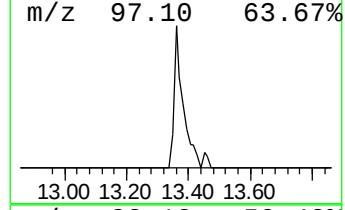
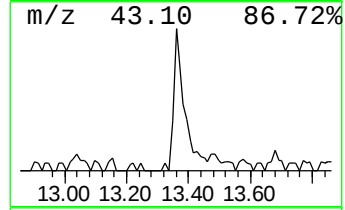
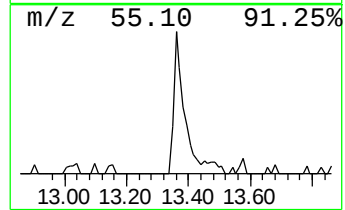
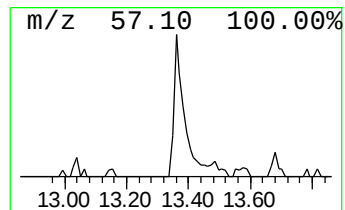
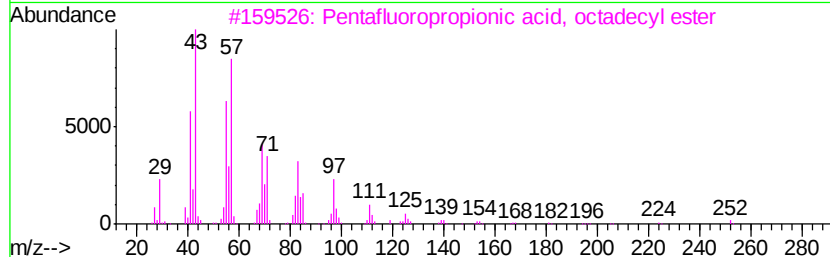
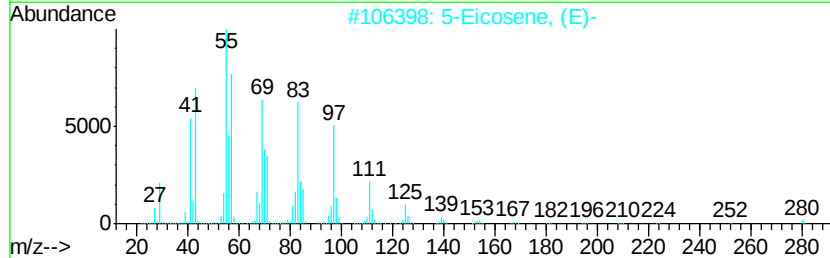
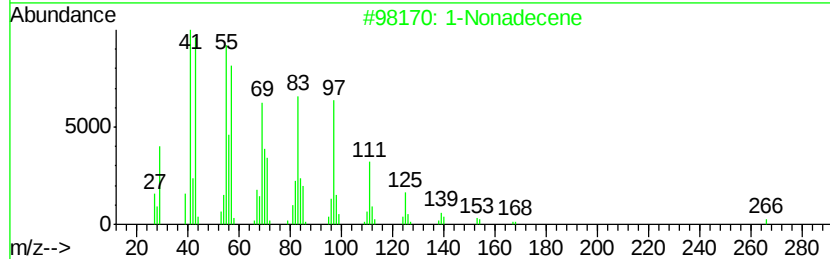
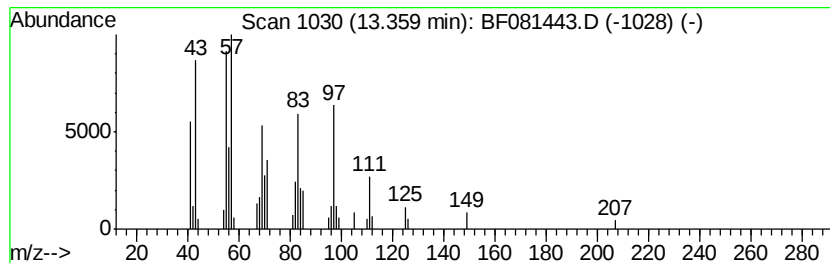
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Peak Number 4 1-Nonadecene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	7.32 ng	128362	Chrysene-d12	13.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Nonadecene	266 C19H38	018435-45-5	91
2	5-Eicosene, (E)-	280 C20H40	074685-30-6	90
3	Pentafluoropropionic acid, octadecyl ester	416 C21H37F5O2	1000280-07-7	83
4	3-Chloropropionic acid, heptadecyl ester	346 C20H39ClO2	1000283-05-1	83
5	Trifluoroacetic acid, n-heptadecyl ester	324 C17H31F3O2	1000216-79-2	83



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.41	73.0	ng	1307480	1	6.35	358327	20.0
unknown6.12	6.12	98.0	ng	1755330	1	6.35	358327	20.0
1-Nonadecene	13.36	7.3	ng	128362	5	13.45	350610	20.0