

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072518\
 Data File : BG035900.D
 Acq On : 26 Jul 2018 8:48
 Operator : JU/SJ
 Sample : J4107-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 071918-DBRI-E-U-S

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_G\METHODS\8270-BG071218.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	5.607	389	395	403	rBV	172781	300133	8.86%	2.210%
2	7.194	657	665	682	rBV	112873	235351	6.95%	1.733%
3	7.558	720	727	741	rBV	360168	662962	19.58%	4.881%
4	8.028	798	807	814	rBV	128514	214652	6.34%	1.580%
5	8.351	853	862	873	rVB	545737	944411	27.89%	6.954%
6	9.185	995	1004	1017	rBV	421640	828783	24.48%	6.102%
7	10.830	1276	1284	1295	rBV	147562	292556	8.64%	2.154%
8	13.274	1692	1700	1710	rBV	1434808	2291012	67.66%	16.869%
9	14.102	1836	1841	1849	rBV	36612	67265	1.99%	0.495%
10	14.648	1928	1934	1942	rBV2	312413	519996	15.36%	3.829%
11	15.935	2149	2153	2158	rBV	41264	54032	1.60%	0.398%
12	16.134	2181	2187	2204	rBV	789992	1338758	39.54%	9.857%
13	17.392	2395	2401	2413	rBV	455466	738981	21.83%	5.441%
14	20.000	2838	2845	2855	rBV	2589036	3385844	100.00%	24.930%
15	21.292	3061	3065	3071	rBV3	46654	82806	2.45%	0.610%
16	21.656	3120	3127	3135	rBV	499238	826863	24.42%	6.088%
17	24.852	3661	3671	3682	rBV3	268781	796874	23.54%	5.867%

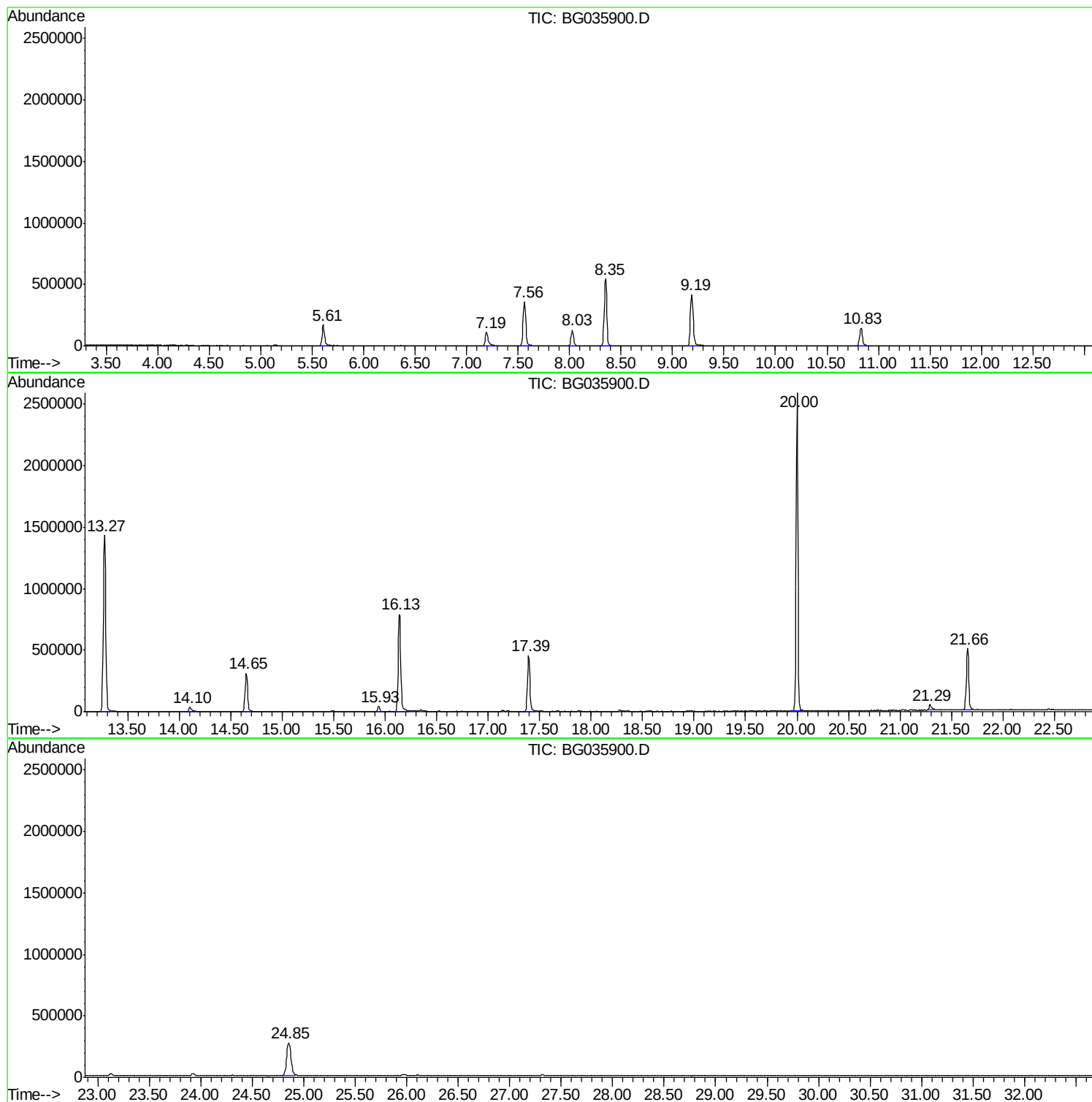
Sum of corrected areas: 13581279

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.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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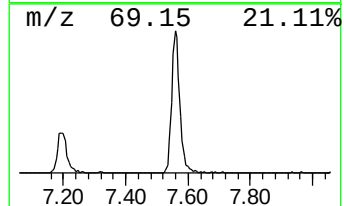
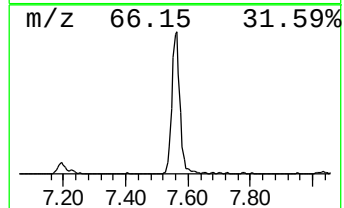
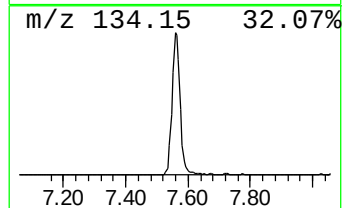
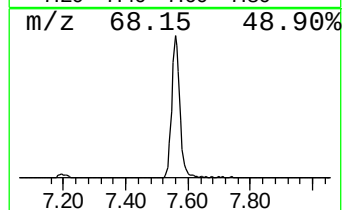
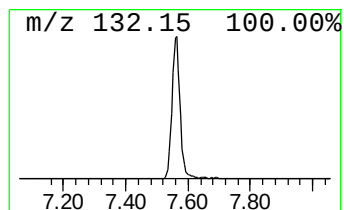
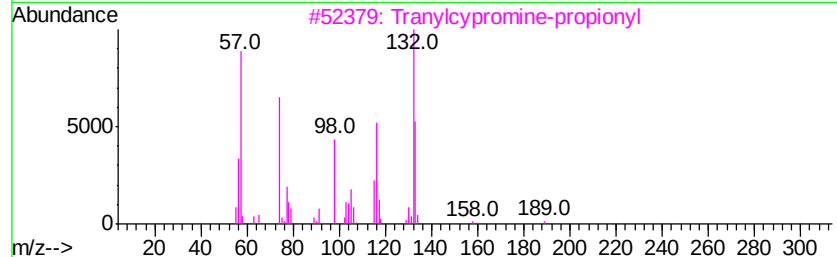
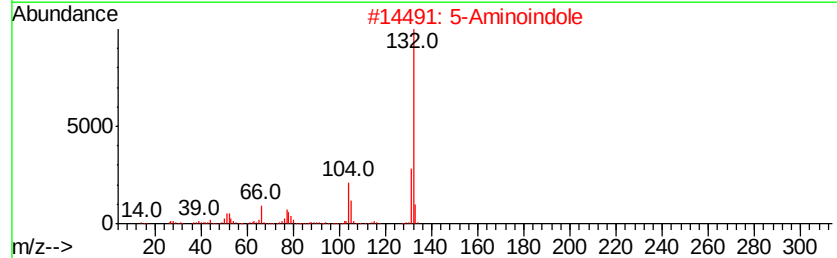
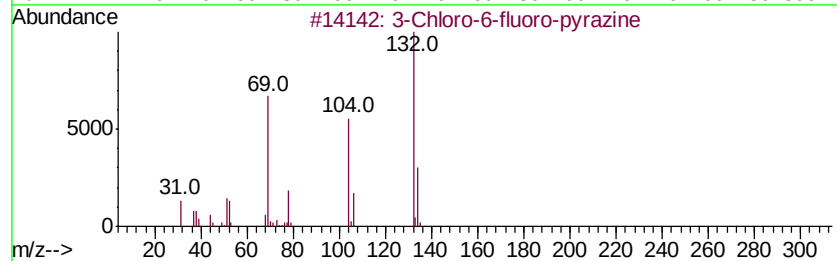
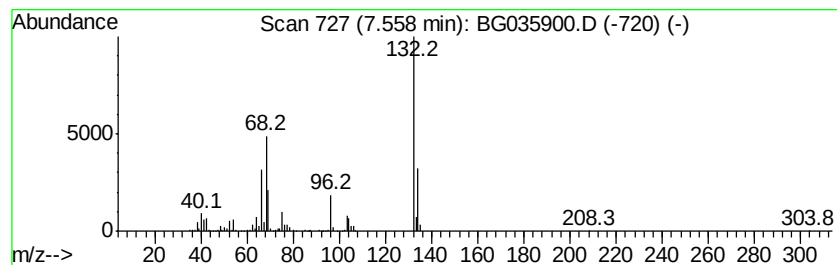
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	61.77 ng	662962	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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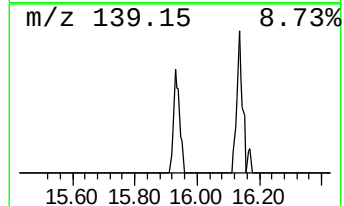
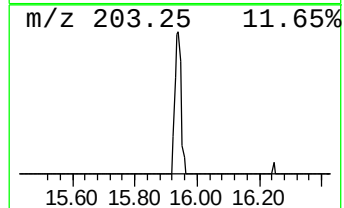
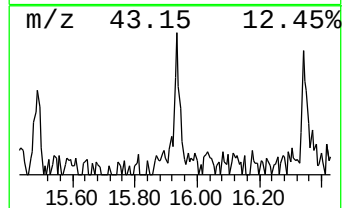
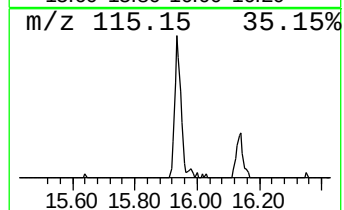
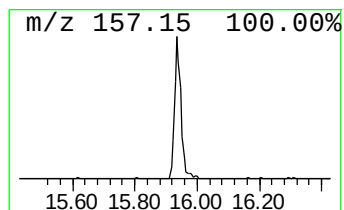
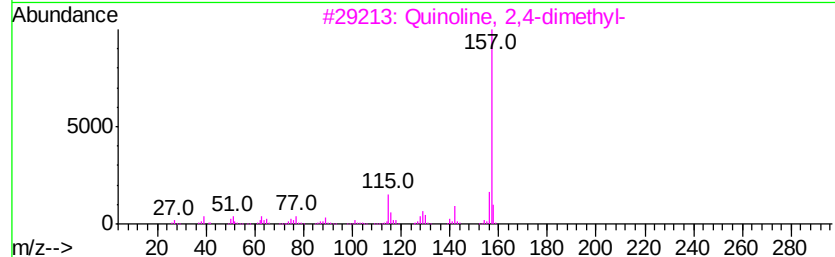
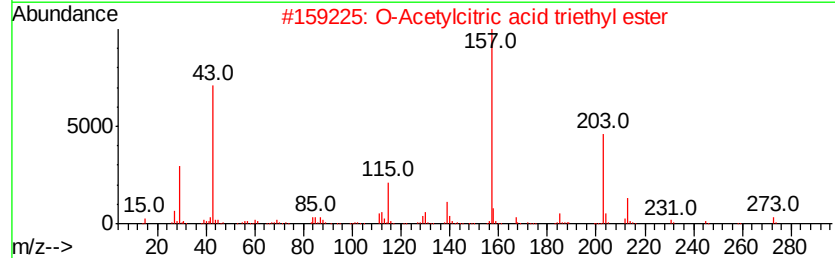
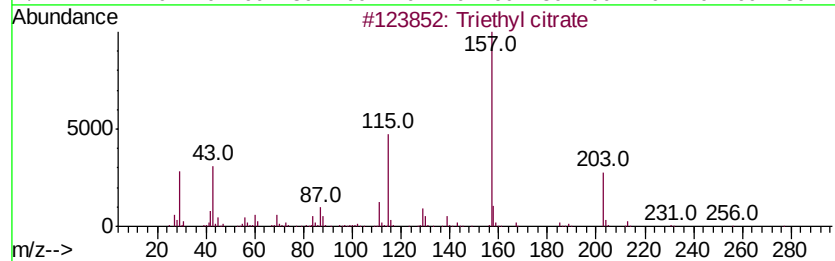
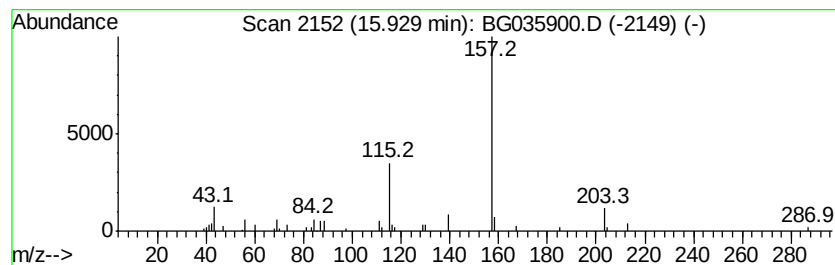
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Peak Number 3 Triethyl citrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	2.08 ng	54032	Acenaphthene-d10	14.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triethyl citrate	276	C12H20O7	000077-93-0	86
2		O-Acetylcitric acid triethyl ester	318	C14H22O8	000077-89-4	64
3		Quinoline, 2,4-dimethyl-	157	C11H11N	001198-37-4	50
4		Piperazine-1-carboxylic acid, 4-...	362	C14H19ClN2O5S	1000303-80-2	40
5		3-Chloro2-fluorobenzoic acid, 4-...	370	C21H32ClF02	1000338-64-7	36



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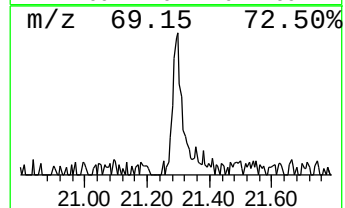
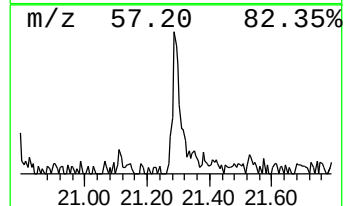
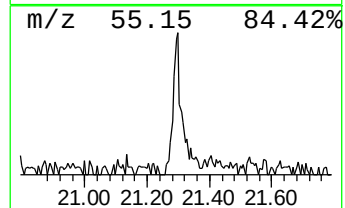
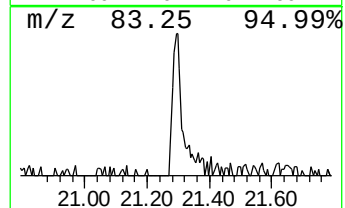
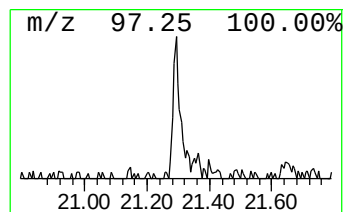
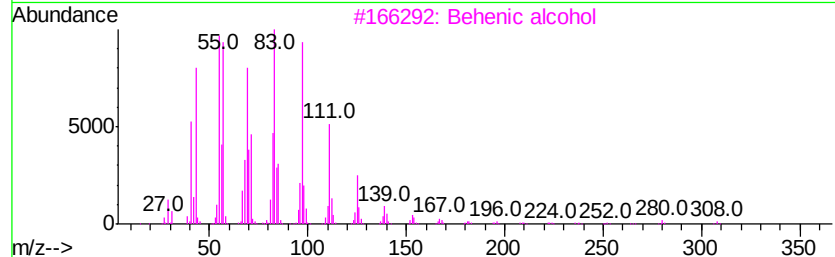
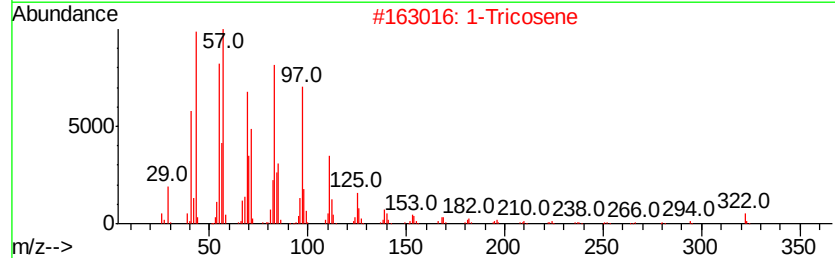
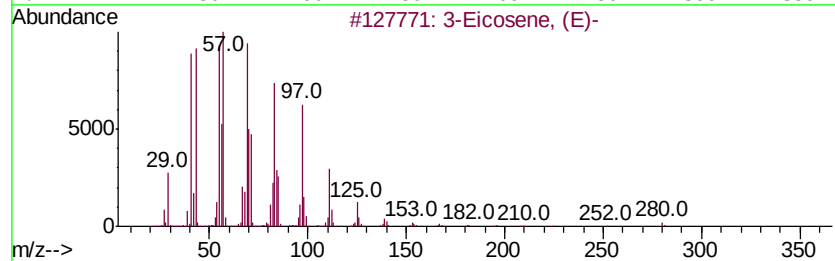
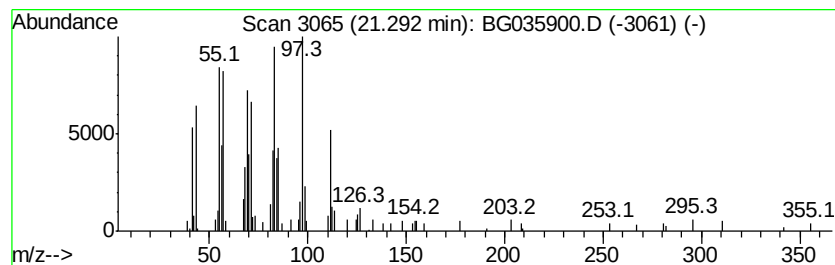
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Peak Number 4 3-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.29	2.00 ng	82806	Chrysene-d12	21.66
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Eicosene, (E)-	280 C20H40	074685-33-9 93
2		1-Tricosene	322 C23H46	018835-32-0 91
3		Behenic alcohol	326 C22H46O	000661-19-8 91
4		Heptafluorobutyric acid, hexadec...	438 C20H33F7O2	006385-15-5 90
5		n-Nonadecanol-1	284 C19H40O	001454-84-8 90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.56	7.56	61.8	ng	662962	1	8.03	214652	20.0
Triethyl citrate	15.93	2.1	ng	54032	3	14.65	519996	20.0
3-Eicosene, (E)-	21.29	2.0	ng	82806	5	21.66	826863	20.0