

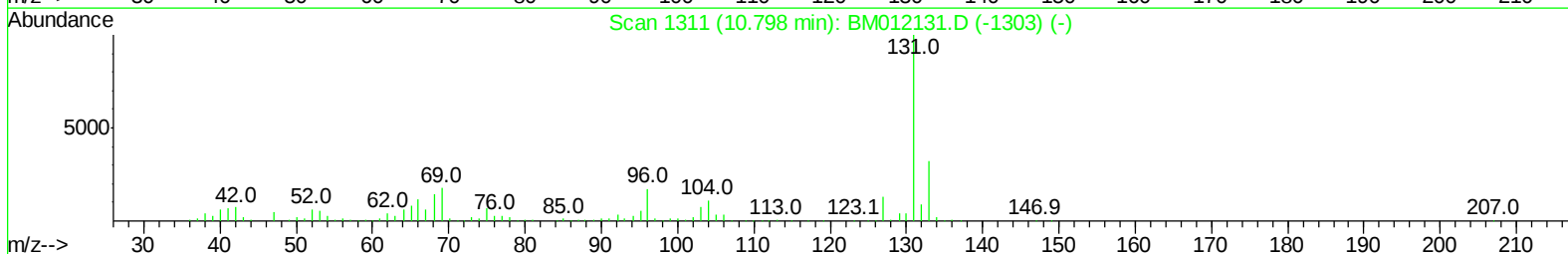
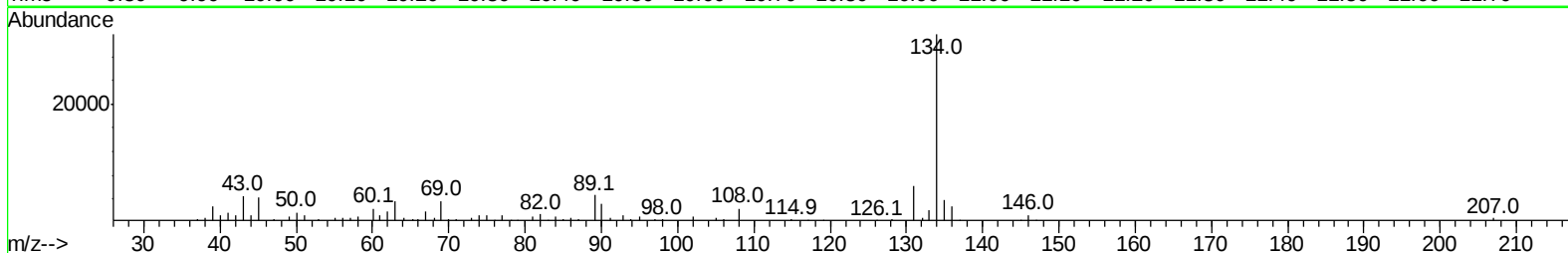
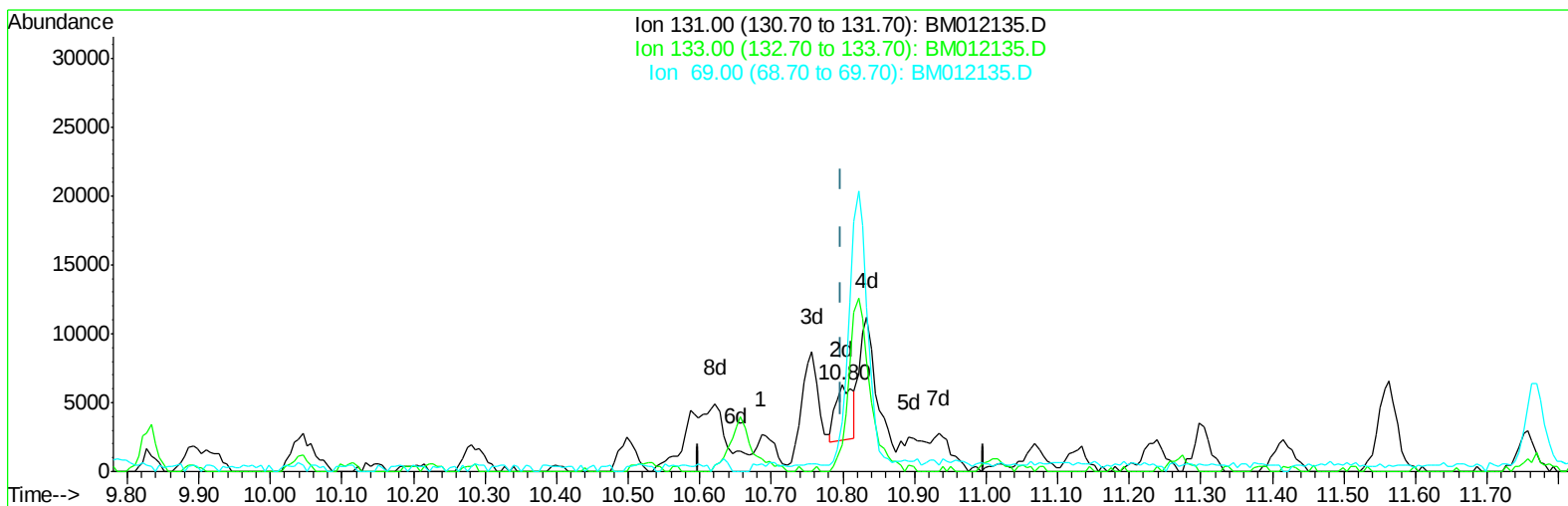
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Data File : BM012135.D
Acq On : 13 Oct 2017 13:54
Operator : SJ/JU
Sample : I5772-03DL 5X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BD9S0DL

Manual Integrations
APPROVED

Sohil
10/15/2017 12:37:32 PM

Quant Time: Oct 14 00:44:00 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 14 00:15:07 2017
Response via : Initial Calibration



TIC: BM012135.D

(29) 4-Chloroaniline-d4 (S)

10.798min (+0.000) 0.49ng/ul m

response 6965

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	31.71
69.00	24.30	58.43#
0.00	0.00	0.00

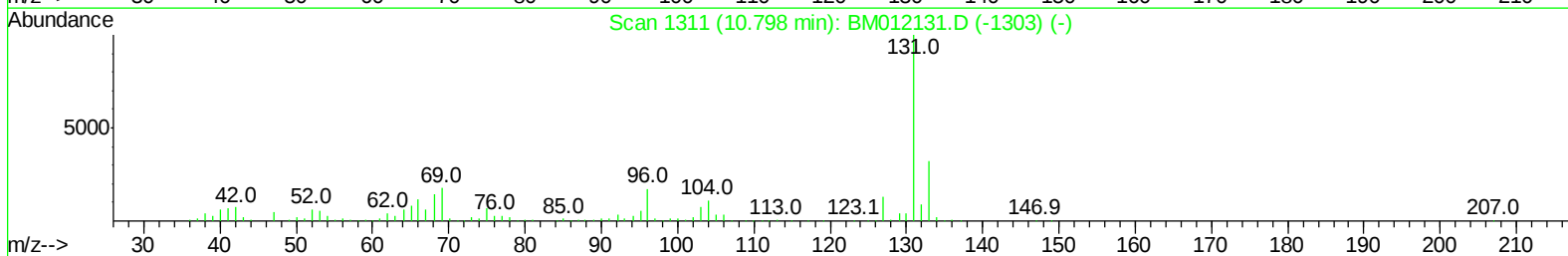
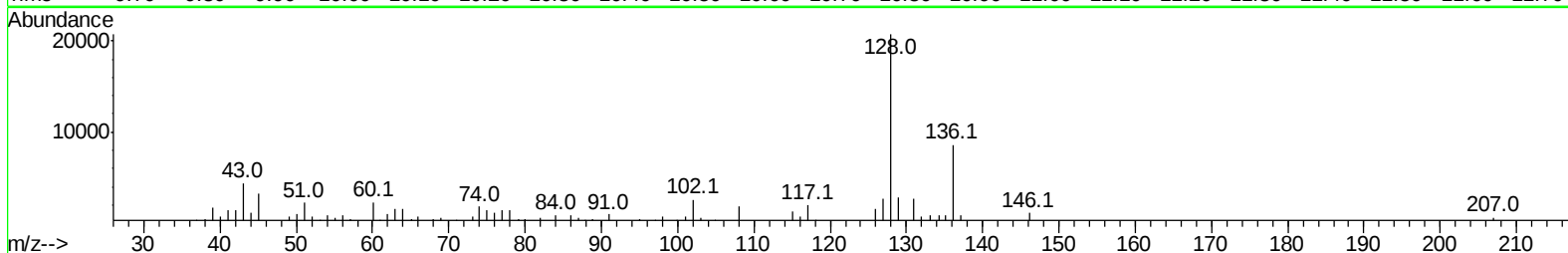
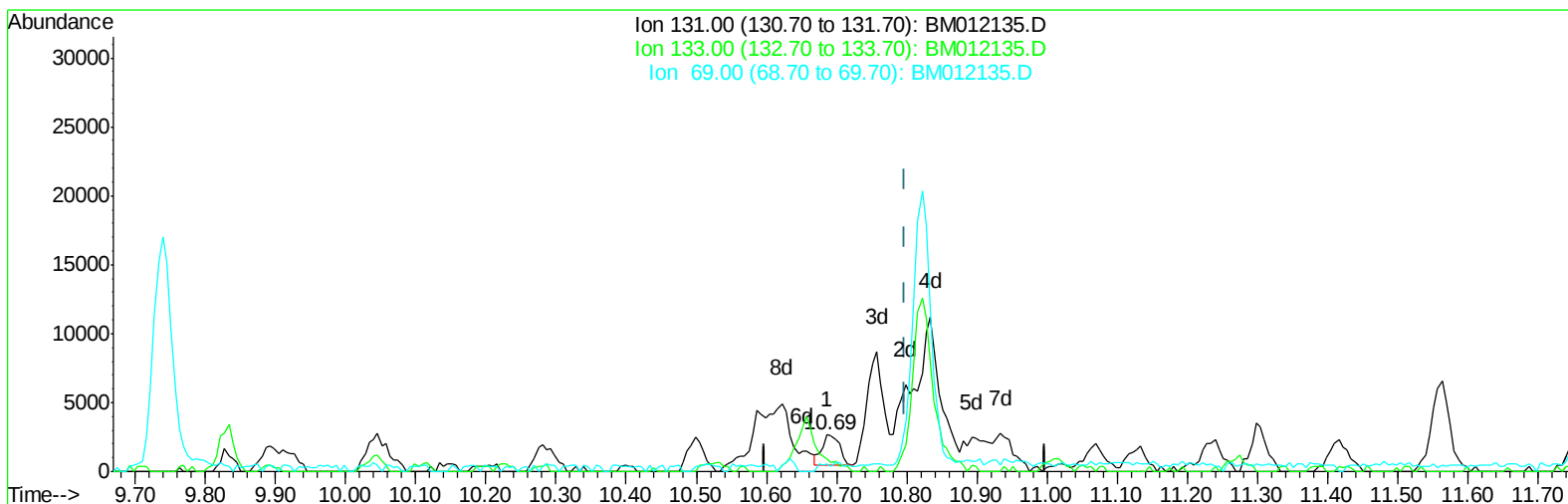
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TIC: BM012135.D

(29) 4-Chloroaniline-d4 (S)
10.687min (-0.112) 0.25ng/ul

response 3594

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	34.96
69.00	24.30	19.99
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	209251	20.00	ng/ul	0.00
18) Naphthalene-d8	10.65	136	886229	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.50	164	530972	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.25	188	1222872	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	1129504	20.00	ng/ul	0.00
83) Perylene-d12	23.76	264	958585	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	2143	0.49	ng/uL	0.00
5) Phenol-d5	7.02	99	29239	1.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.17	67	70157	6.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.38	132	84577	5.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.57	113	54312	3.72	ng/ul	0.00
19) Nitrobenzene-d5	9.02	128	51857	7.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.74	143	55843	7.07	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.29	165	101496	6.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.80	131	6965m	0.49	ng/ul	0.00
43) Dimethylphthalate-d6	13.90	166	364109	7.76	ng/ul	0.00
46) Acenaphthylene-d8	14.19	160	440714	7.76	ng/ul	0.00
51) 4-Nitrophenol-d4	14.75	143	7626	1.08	ng/ul	0.02
57) Fluorene-d10	15.49	176	304511	7.91	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.65	200	8016	0.96	ng/ul	0.00
70) Anthracene-d10	17.34	188	491082	8.12	ng/ul	0.00
76) Pyrene-d10	19.64	212	528895	9.22	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.62	264	364624	7.89	ng/ul	0.00

Target Compounds

					Ovalue
28) Naphthalene	10.70	128	91013	1.985	ng/ul 98
40) 1,1'-Biphenyl	13.33	154	684084	15.266	ng/ul 99
49) Acenaphthene	14.56	153	996375	26.518	ng/ul 100
53) Dibenzofuran	14.90	168	1666861	30.809	ng/ul 98
58) Fluorene	15.55	166	240130	5.344	ng/ul# 99
69) Phenanthrene	17.29	178	582474	8.373	ng/ul 98
71) Anthracene	17.38	178	98899	1.373	ng/ul 97
72) Carbazole	17.66	167	1083678	17.779	ng/ul 99
73) Di-n-butylphthalate	18.19	149	177511	2.098	ng/ul 98
74) Fluoranthene	19.31	202	134670	1.604	ng/ul# 93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

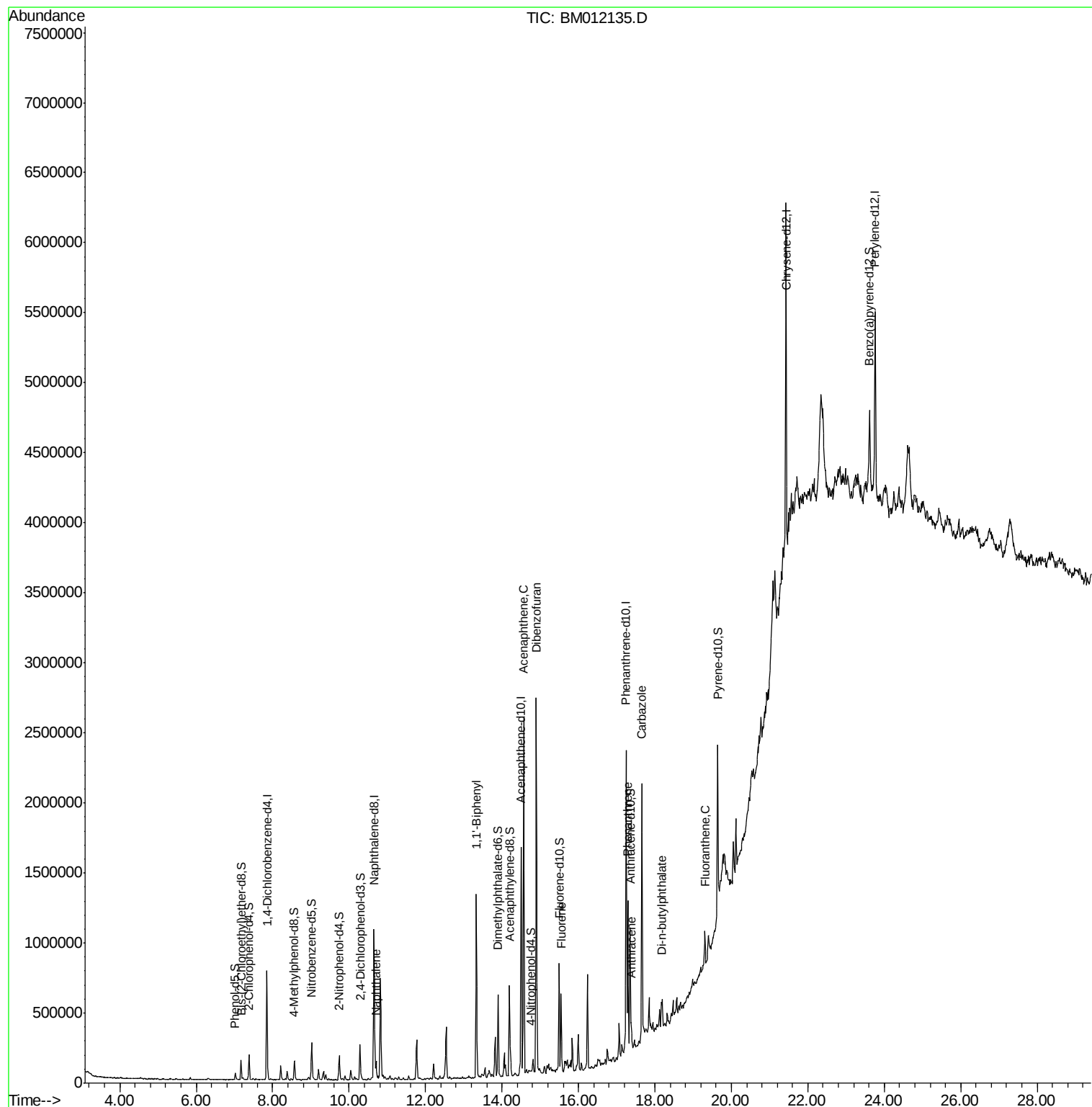
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Manual Integrations
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