

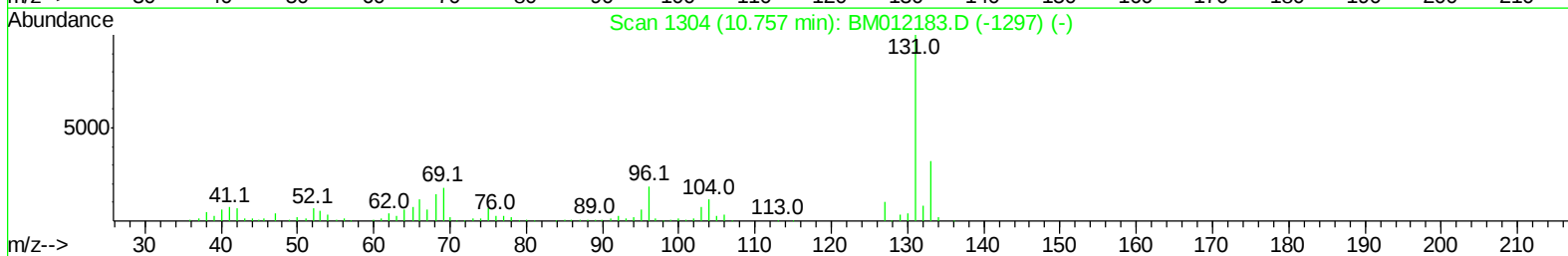
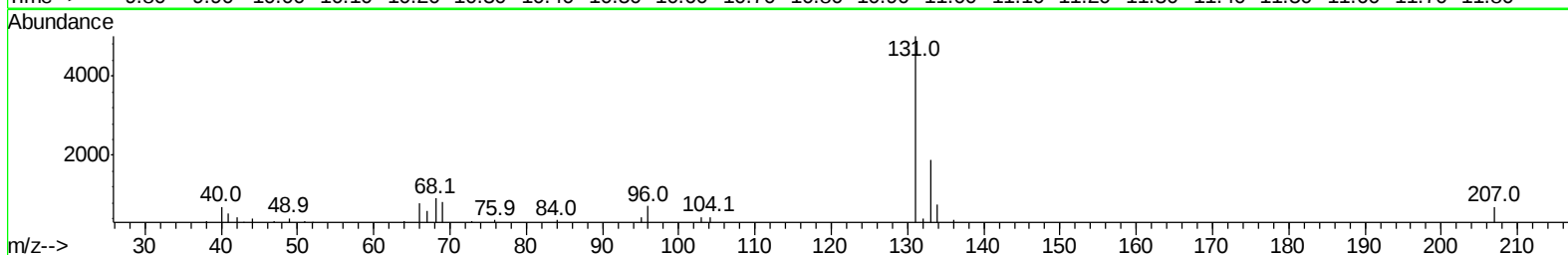
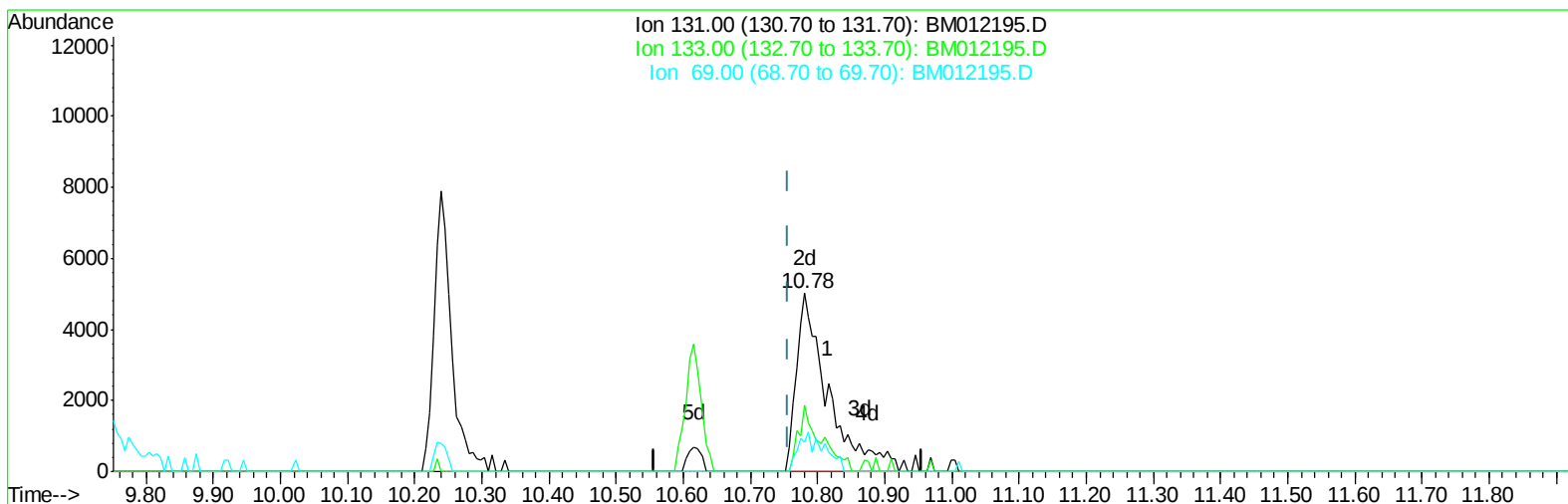
Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM101417\
Data File : BM012195.D
Acq On : 15 Oct 2017 06:08
Operator : SJ/JU
Sample : I5840-08
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AF0

Manual Integrations
APPROVED

Sohil
10/16/2017 7:48:37 PM

Quant Time: Oct 16 01:32:27 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Oct 15 06:19:21 2017
Response via : Initial Calibration



TIC: BM012195.D

(29) 4-Chloroaniline-d4 (S)

10.781min (+0.024) 1.30ng/ul m

response 16463

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	37.42
69.00	24.30	16.43#
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	7.81	152	171665	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	797459	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.46	164	517251	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.21	188	1231606	20.00	ng/ul	0.00
75) Chrysene-d12	21.39	240	1200880	20.00	ng/ul	0.00
83) Perylene-d12	23.70	264	1005377	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.22	96	16311	4.52	ng/uL	0.00
5) Phenol-d5	6.98	99	95704	6.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	233481	27.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	301910	25.55	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	204459	17.05	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	189356	31.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	235066	33.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	416429	30.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.78	131	16463m	1.30	ng/ul	0.02
43) Dimethylphthalate-d6	13.87	166	1526296	33.39	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	1859893	33.61	ng/ul	0.00
51) 4-Nitrophenol-d4	14.70	143	30636	4.46	ng/ul	0.02
57) Fluorene-d10	15.46	176	1306686	34.82	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.60	200	159438	18.86	ng/ul	0.00
70) Anthracene-d10	17.31	188	2076250	34.09	ng/ul	0.00
76) Pyrene-d10	19.60	212	2416067	39.63	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.55	264	1695369	34.98	ng/ul	0.00

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

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