

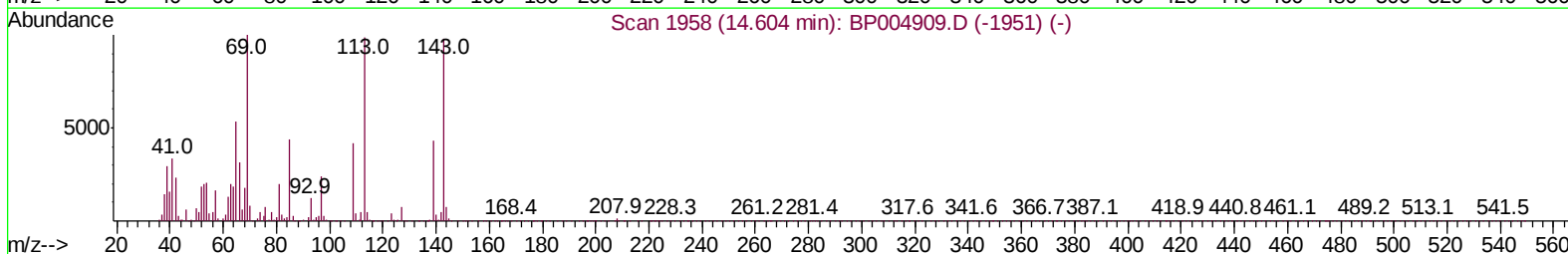
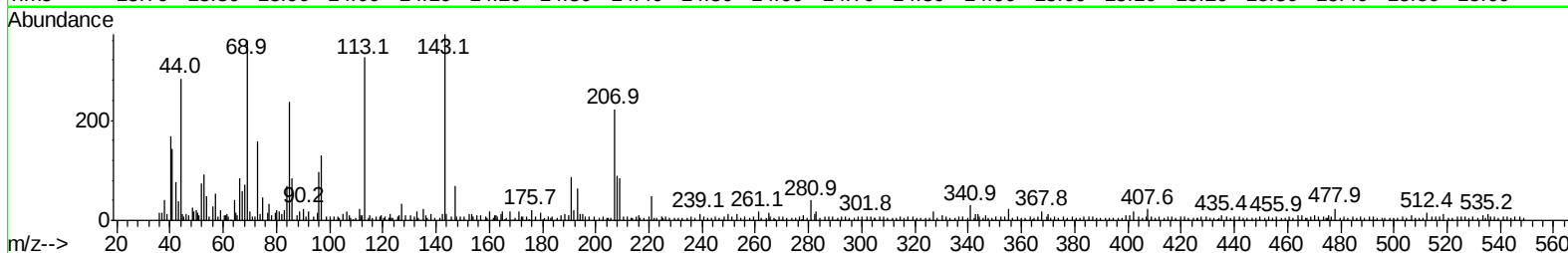
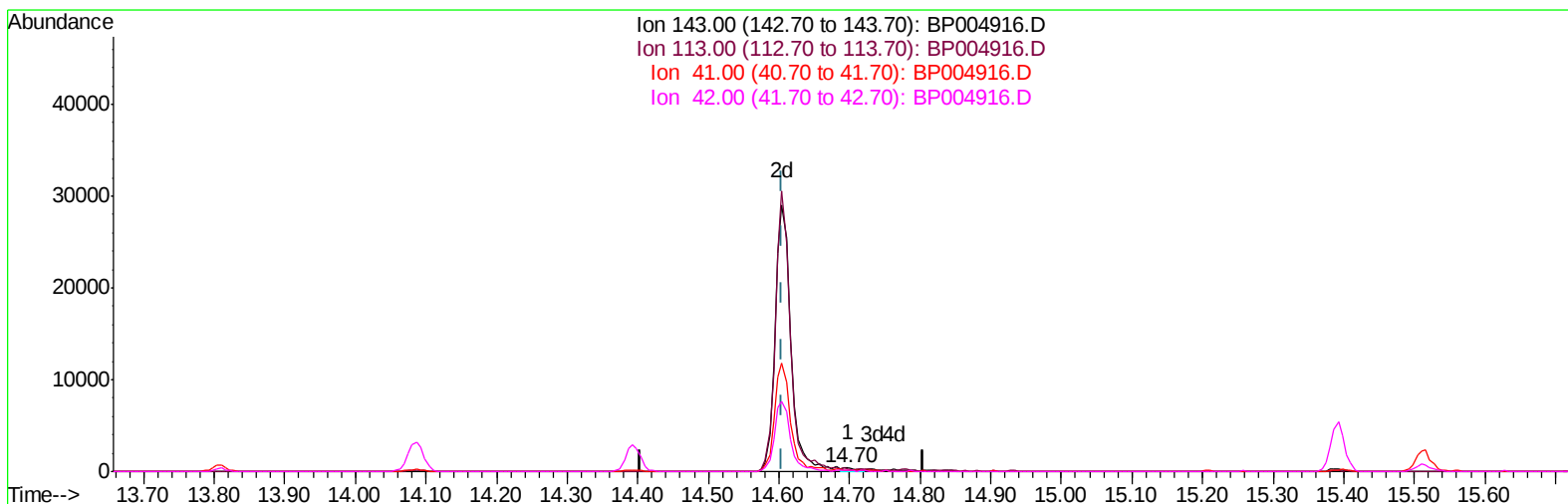
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030521\
Data File : BP004916.D
Acq On : 05 Mar 2021 15:02
Operator : CG/JU
Sample : PB134929BL
Misc : SOM-MDL-W-BL-04
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK29

Manual Integrations
APPROVED

mohammad
3/8/2021 9:36:34 AM

Quant Time: Mar 05 15:31:36 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030521MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 05 13:02:07 2021
Response via : Initial Calibration



TIC: BP004916.D

(52) 4-Nitrophenol-d4 (S)

14.698min (+0.094) 0.36ng/ul

response 418

Ion	Exp%	Act%
143.00	100	100
113.00	100.70	87.37
41.00	34.00	38.17
42.00	24.20	20.70

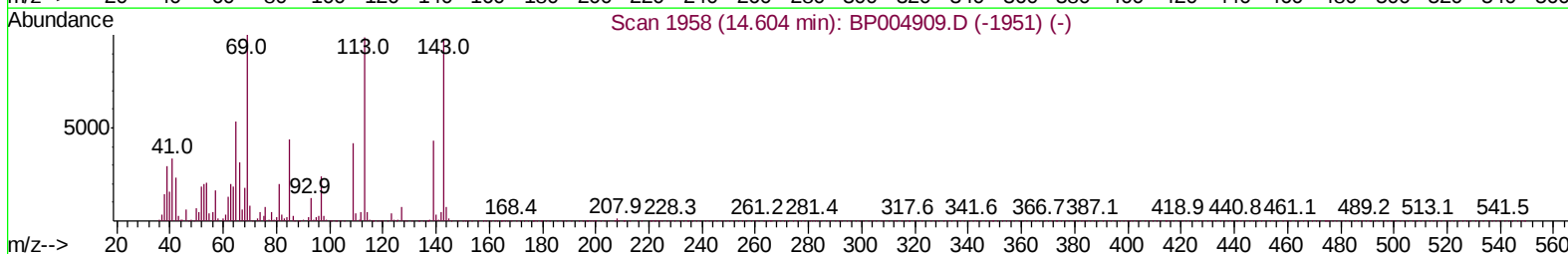
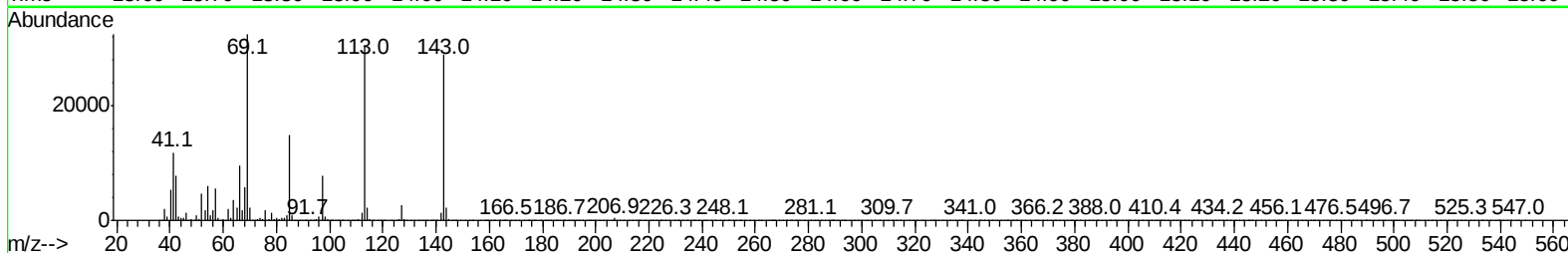
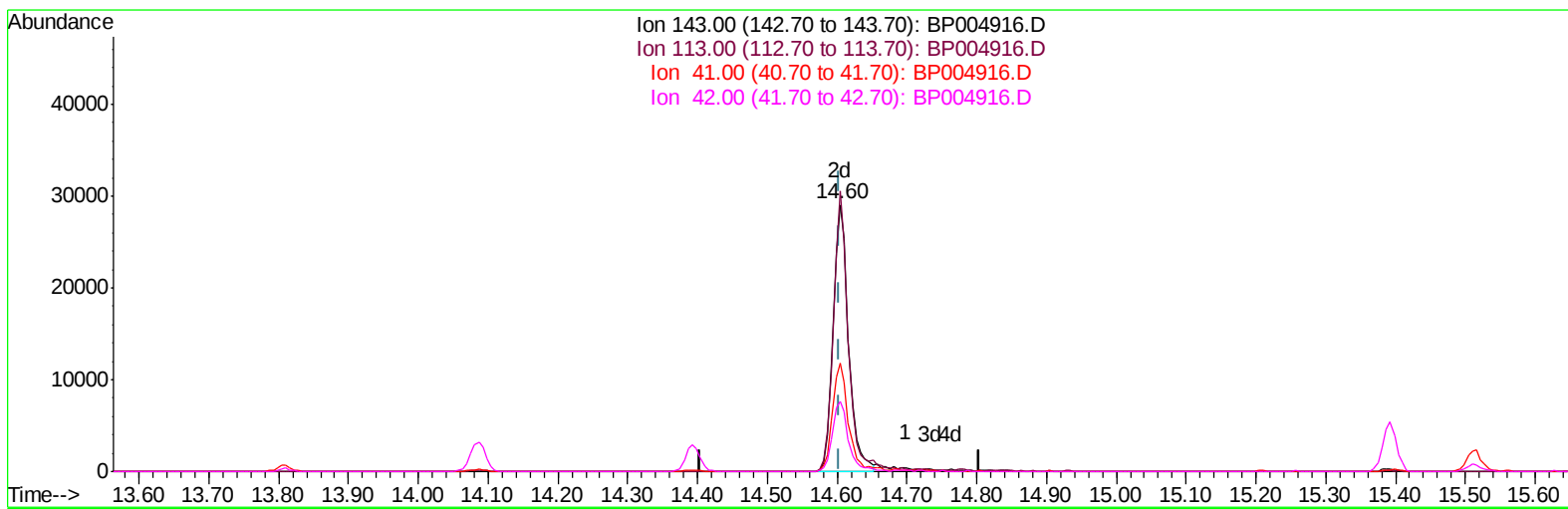
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TIC: BP004916.D

(52) 4-Nitrophenol-d4 (S)

14.604min (-0.000) 37.63ng/ul m

response 44206

Ion	Exp%	Act%
143.00	100	100
113.00	100.70	105.24
41.00	34.00	40.88#
42.00	24.20	26.51

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	35477	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	139381	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	93006	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.15	188	207996	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	230792	20.00	ng/ul	0.00
86) Perylene-d12	23.54	264	237962	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	6272	7.18	ng/uL	0.00
5) Phenol-d5	6.92	99	98981	36.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	53690	38.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.28	132	78202	36.16	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	78984	36.04	ng/ul	0.00
19) Nitrobenzene-d5	8.90	128	37699	37.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	42992	37.61	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.16	165	76985	34.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.68	131	106719	43.46	ng/ul	0.00
44) Dimethylphthalate-d6	13.81	166	285352	42.65	ng/ul	0.00
47) Acenaphthylene-d8	14.09	160	314149	38.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	44206m	37.63	ng/ul	0.00
58) Fluorene-d10	15.39	176	232844	39.05	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	48752	35.80	ng/ul	0.00
71) Anthracene-d10	17.25	188	397261	42.54	ng/ul	0.00
79) Pyrene-d10	19.50	212	464978	44.69	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.40	264	513970	42.50	ng/ul	0.00

Target Compounds Ovalue

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

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