

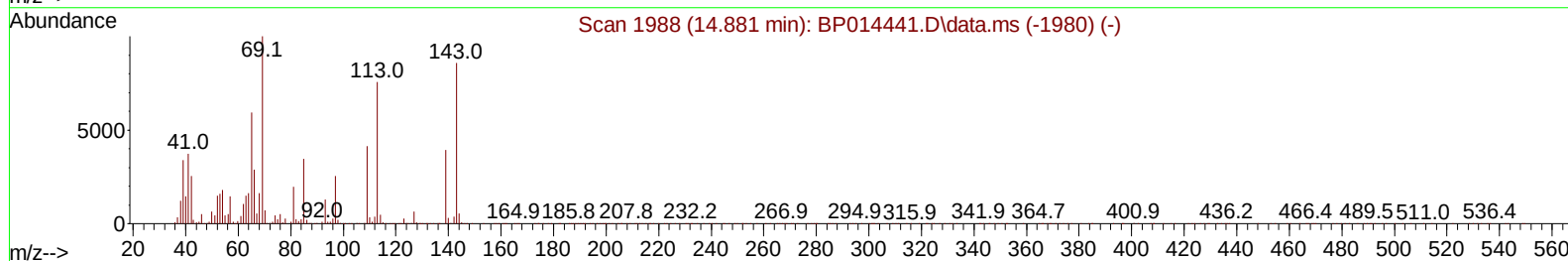
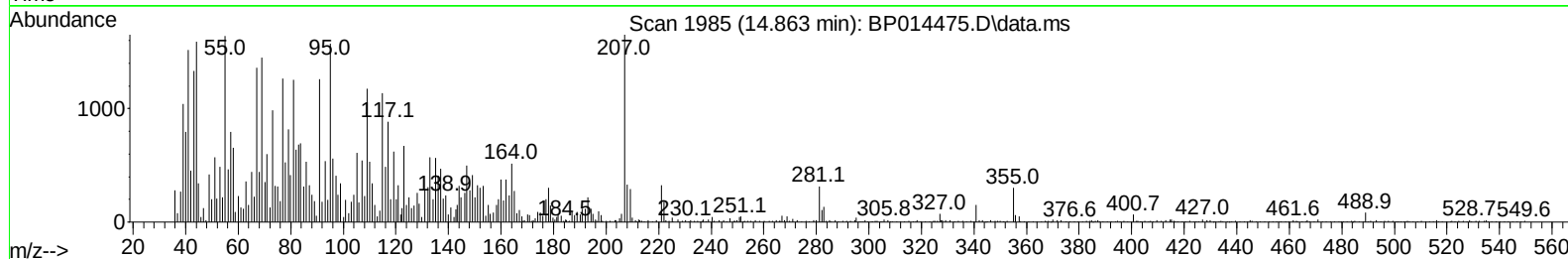
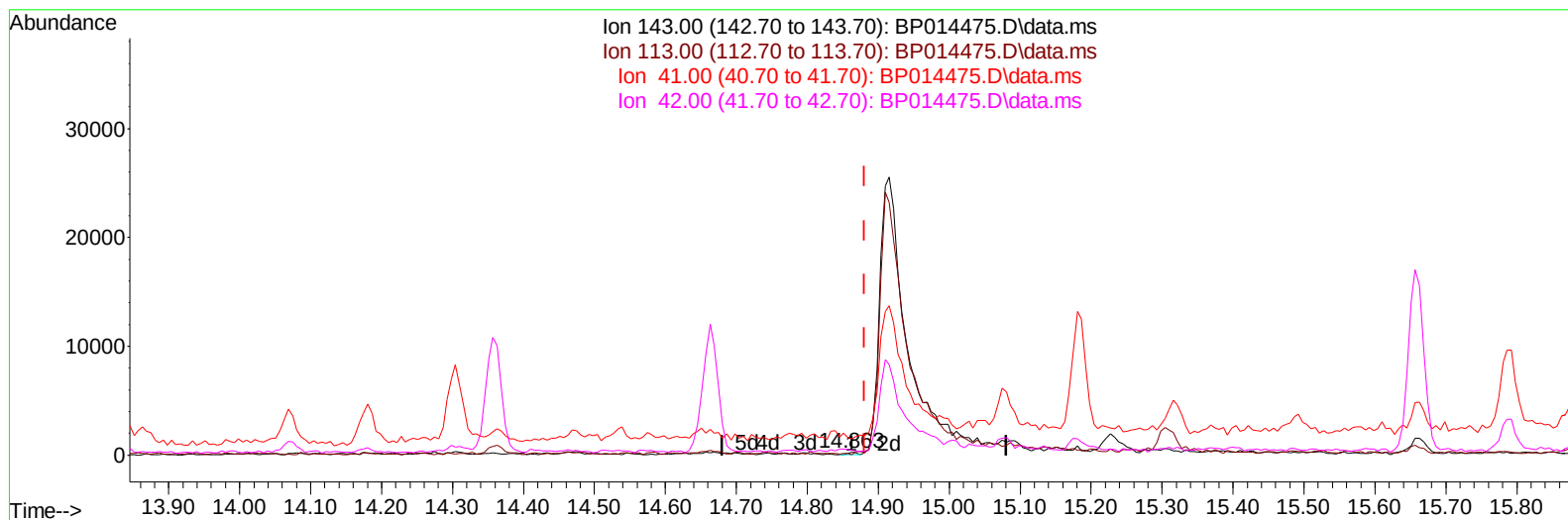
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014475.D
Acq On : 01 Apr 2023 09:25
Operator : CG/JU
Sample : 02158-06
Misc :
ALS Vial : 85 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-5(20230329)

Manual IntegrationsAPPROVED

Quant Time: Apr 03 00:44:15 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014475.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (-0.018) 0.04 ng/u1

response 223

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	19.85#
41.00	44.20	569.29#
42.00	31.50	168.91#

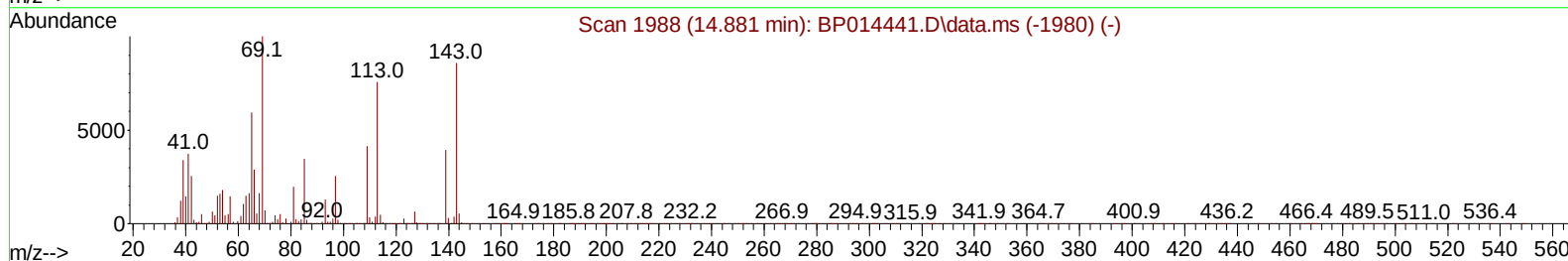
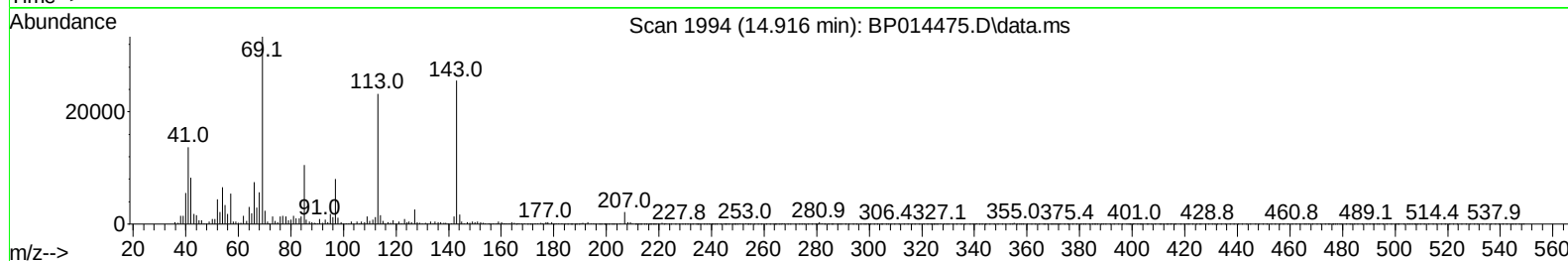
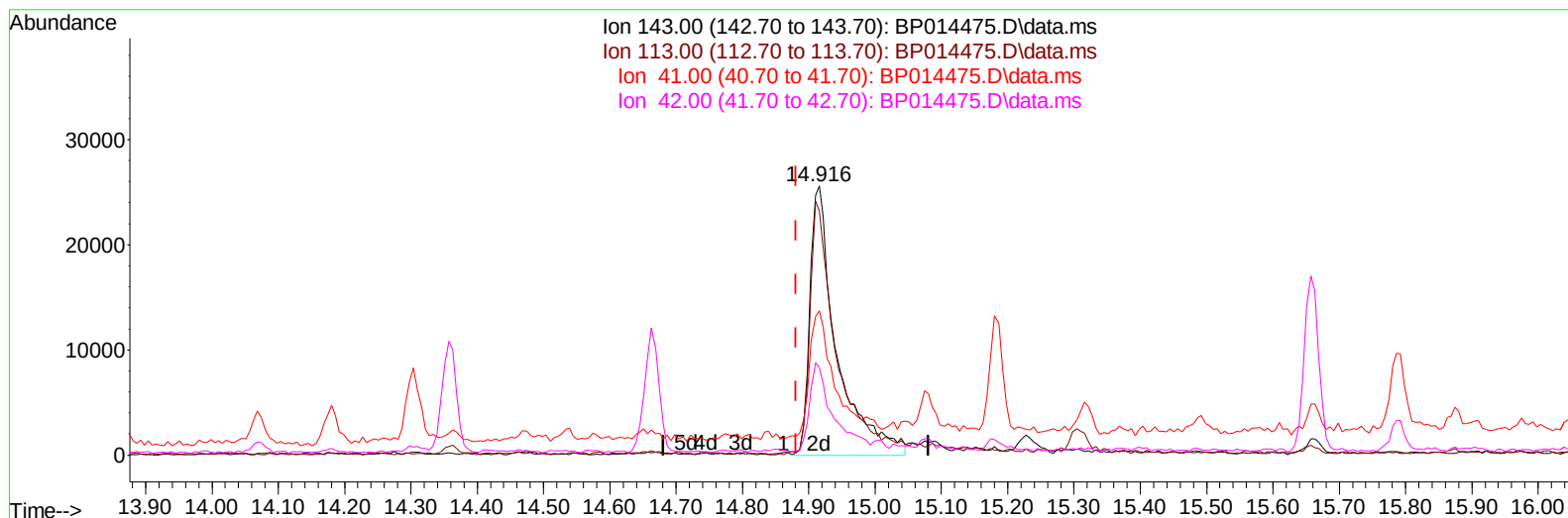
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(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.035) 11.60 ng/ul m

response 71438

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	90.77
41.00	44.20	53.70#
42.00	31.50	32.54

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	170702	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	693795	20.000	ng/ul	0.03
38) Acenaphthene-d10	14.663	164	502826	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.427	188	1101555	20.000	ng/ul	0.00
79) Chrysene-d12	21.515	240	820999	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	771544	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	20683	4.699	ng/uL	0.00
4) Pyridine-d5	3.828	84	129355	11.467	ng/ul	0.01
7) Phenol-d5	7.187	99	201816	12.819	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	245919	24.347	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	250412	22.007	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	221234	17.422	ng/ul	0.01
21) Nitrobenzene-d5	9.193	128	151348	27.022	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	174224	27.166	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	302784	25.923	ng/ul	0.03
31) 4-Chloroaniline-d4	10.992	131	301751	19.613	ng/ul	0.01
46) Dimethylphthalate-d6	14.069	166	1074591	26.357	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1178884	26.059	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	71438m	11.598	ng/ul	0.03
60) Fluorene-d10	15.657	176	899534	26.457	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	146171	18.034	ng/ul	0.00
73) Anthracene-d10	17.527	188	1425624	26.742	ng/ul	0.00
81) Pyrene-d10	19.757	212	1546858	28.050	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1090449	26.602	ng/ul	0.00

Target Compounds Qvalue

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

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