

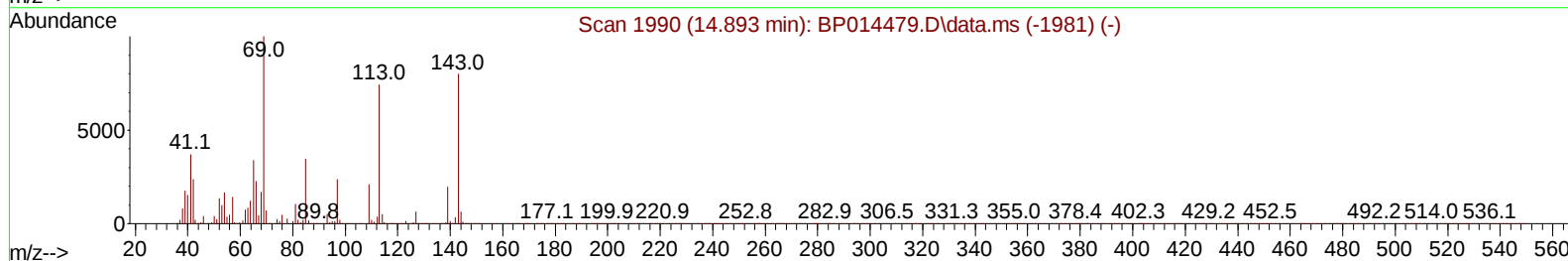
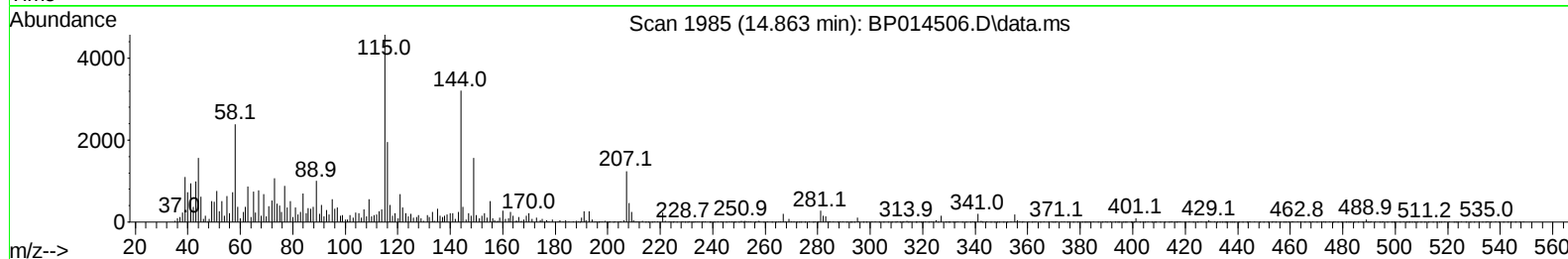
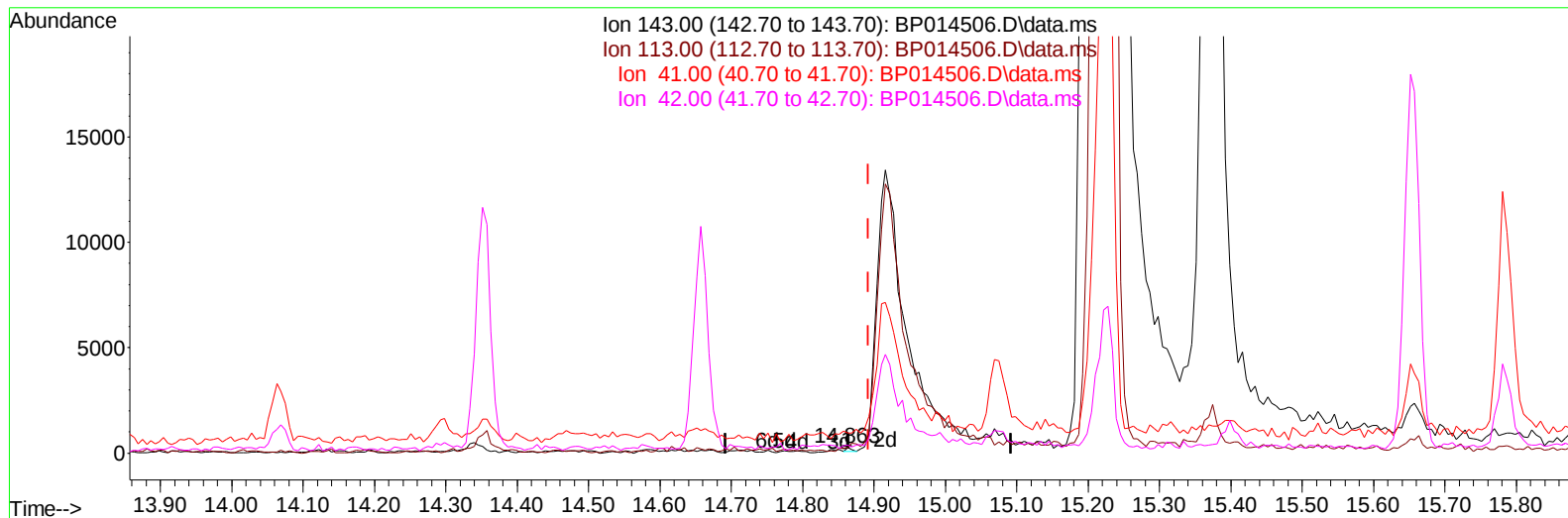
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014506.D
Acq On : 04 Apr 2023 01:58
Operator : CG/JU
Sample : 02158-12
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-9S(20230330)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 02:36:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 04 01:34:36 2023
Response via : Initial Calibration

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



TIC: BP014506.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (-0.029) 0.02 ng/u1

response 84

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	105.24
41.00	44.20	383.06#
42.00	31.50	147.98#

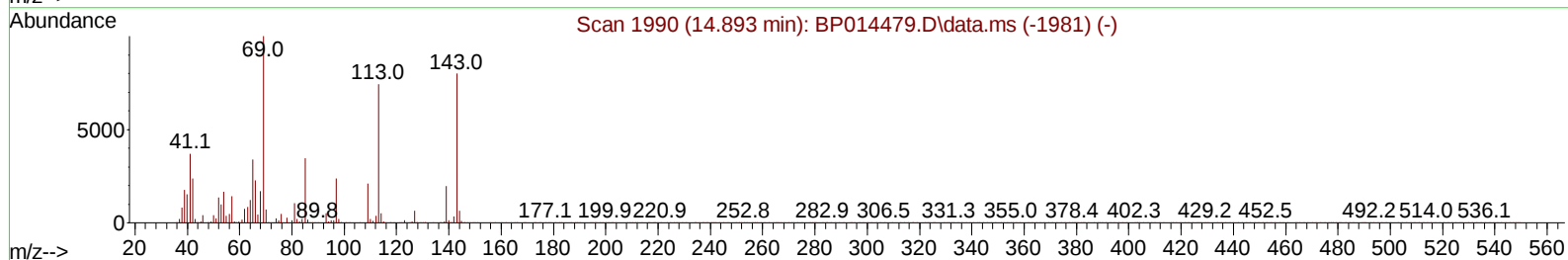
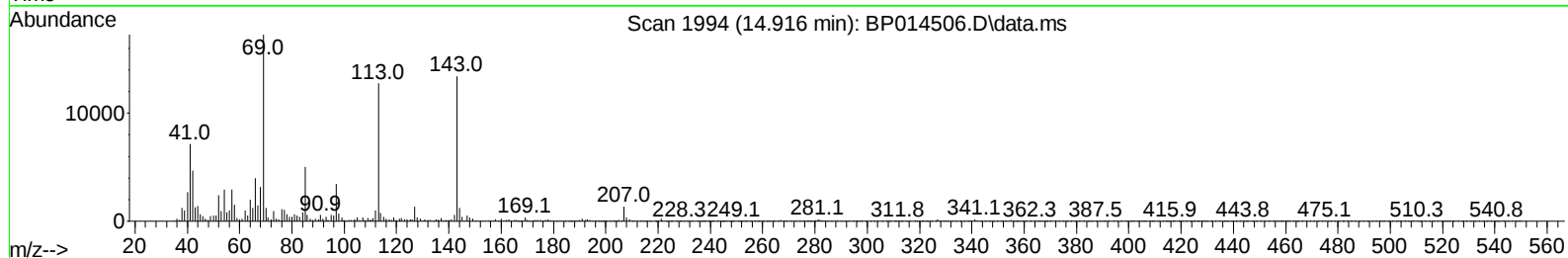
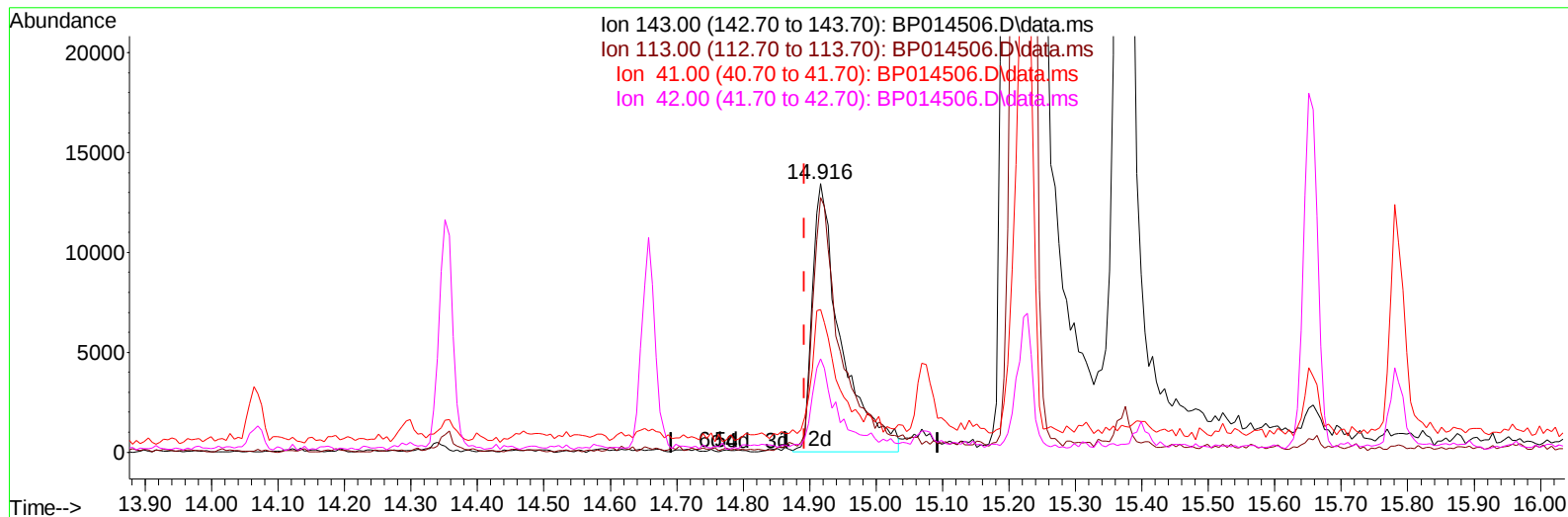
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(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.024) 8.30 ng/ul m

response 40327

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.93
41.00	44.20	53.09#
42.00	31.50	34.82

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	133061	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	583266	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	396467	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	916966	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	1005564	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	1019859	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	14875	4.335	ng/uL	0.00
4) Pyridine-d5	3.828	84	77391	8.801	ng/ul	0.01
7) Phenol-d5	7.175	99	119283	9.720	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	258682	32.855	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	248874	28.060	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	222532	22.481	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	150135	31.885	ng/ul	0.00
24) 2-Nitrophenol-d4	9.911	143	172849	32.058	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.458	165	300595	30.613	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	298314	23.064	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	1077880	33.530	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	1152969	32.324	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	40327m	8.304	ng/ul	0.02
60) Fluorene-d10	15.657	176	891267	33.247	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	181521	26.903	ng/ul	0.00
73) Anthracene-d10	17.522	188	1523721	34.336	ng/ul	0.00
81) Pyrene-d10	19.751	212	1963810	29.075	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.863	264	1878898	34.676	ng/ul	0.00

Target Compounds Qvalue

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

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