

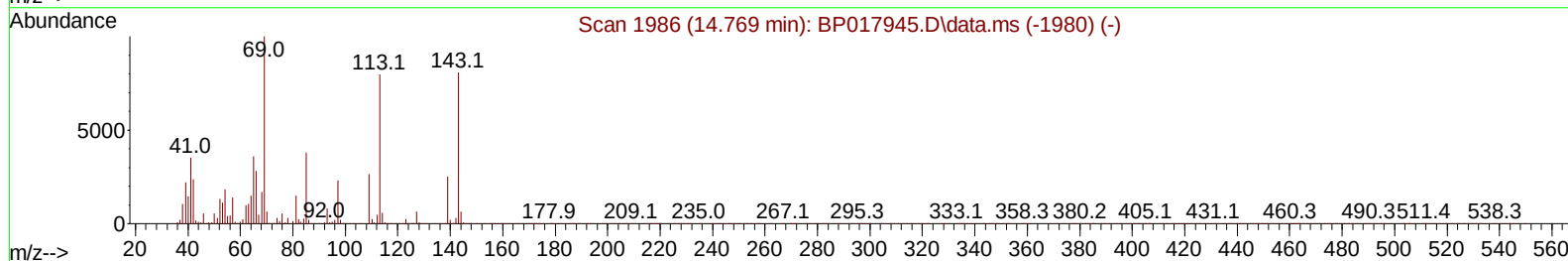
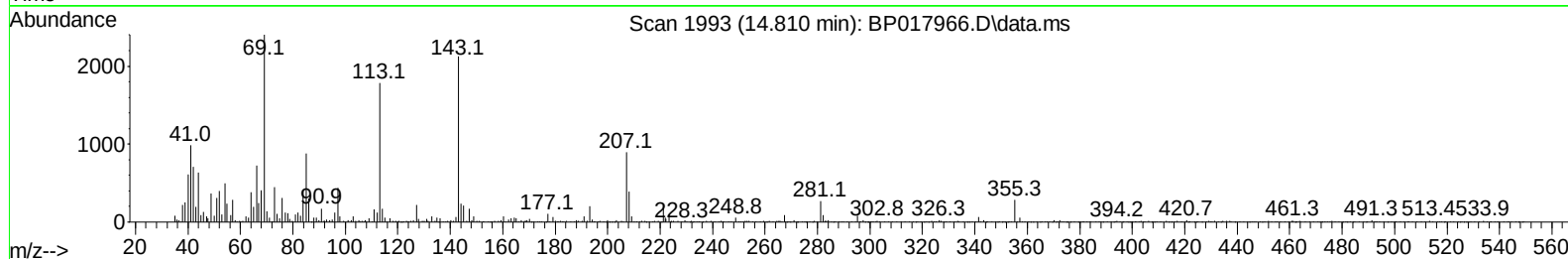
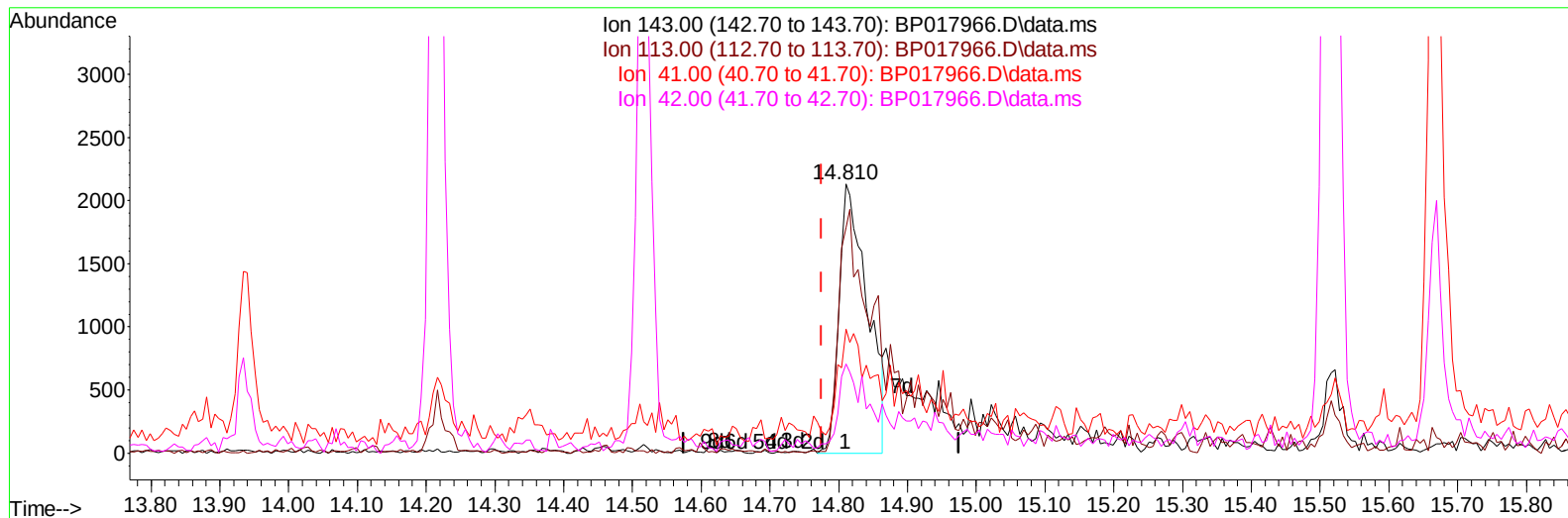
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
Data File : BP017966.D
Acq On : 08 Nov 2023 17:26
Operator : MA/JU
Sample : 05013-13
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4998

Manual IntegrationsAPPROVED

Quant Time: Nov 08 21:43:50 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 07 03:40:56 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/09/2023
Supervised By :mohammad ahmed 11/11/2023



TIC: BP017966.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.035) 2.89 ng/u1

response 6032

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	83.80
41.00	43.00	46.10
42.00	28.20	33.19

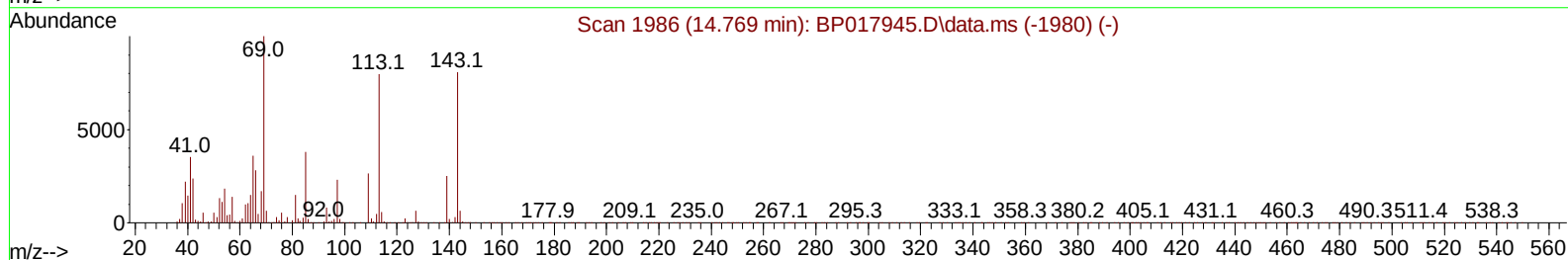
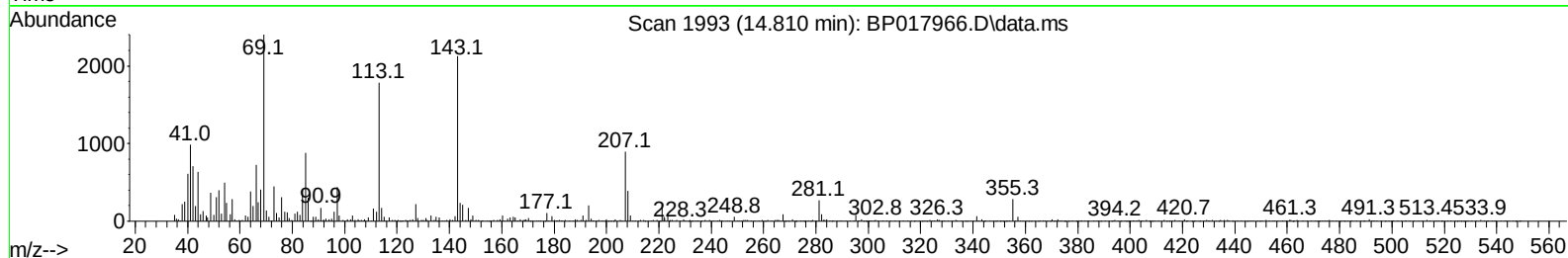
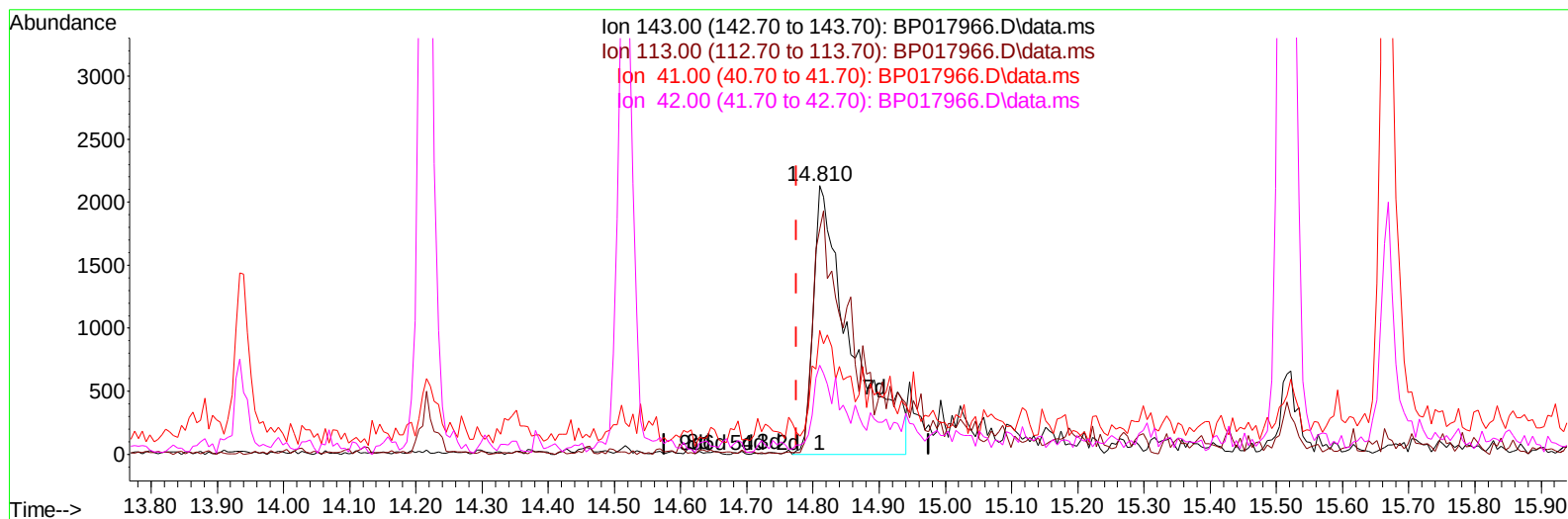
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(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.035) 4.00 ng/ul m

response 8356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	83.80
41.00	43.00	46.10
42.00	28.20	33.19

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Quant Time: Nov 08 22:34:11 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 07 03:40:56 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	57490	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	243410	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	154796	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	345691	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	329267	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	367699	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	6576	4.378	ng/uL	0.00
4) Pyridine-d5	3.687	84	26067	6.253	ng/ul	0.01
7) Phenol-d5	7.028	99	29593	5.489	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	114243	35.129	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	107268	25.894	ng/ul	-0.01
15) 4-Methylphenol-d8	8.575	113	66601	14.946	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	72680	37.364	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	77824	34.736	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	132970	32.374	ng/ul	0.00
31) 4-Chloroaniline-d4	10.852	131	144723	24.315	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	552743	40.723	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	557612	37.439	ng/ul	0.00
54) 4-Nitrophenol-d4	14.810	143	8356m	4.004	ng/ul	0.04
60) Fluorene-d10	15.516	176	447920	39.427	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	80113	31.690	ng/ul	0.00
73) Anthracene-d10	17.392	188	726828	40.438	ng/ul	0.00
81) Pyrene-d10	19.639	212	923141	44.165	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	882578	42.958	ng/ul	0.00

Target Compounds Qvalue

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

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