

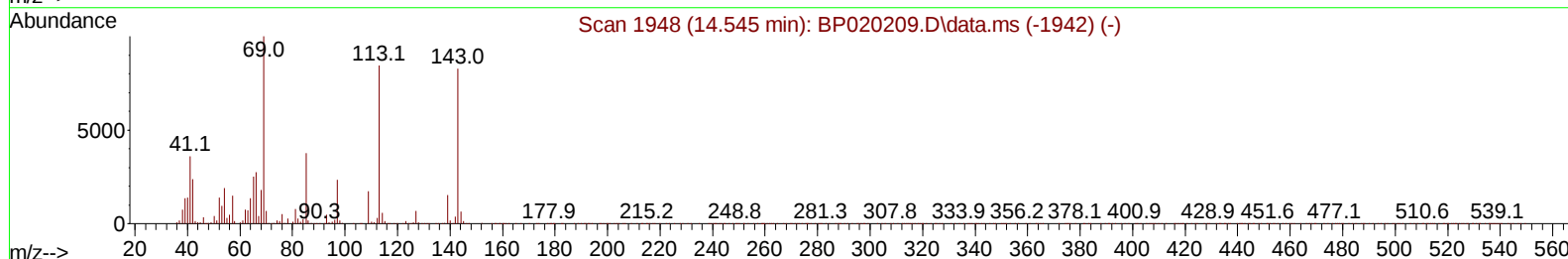
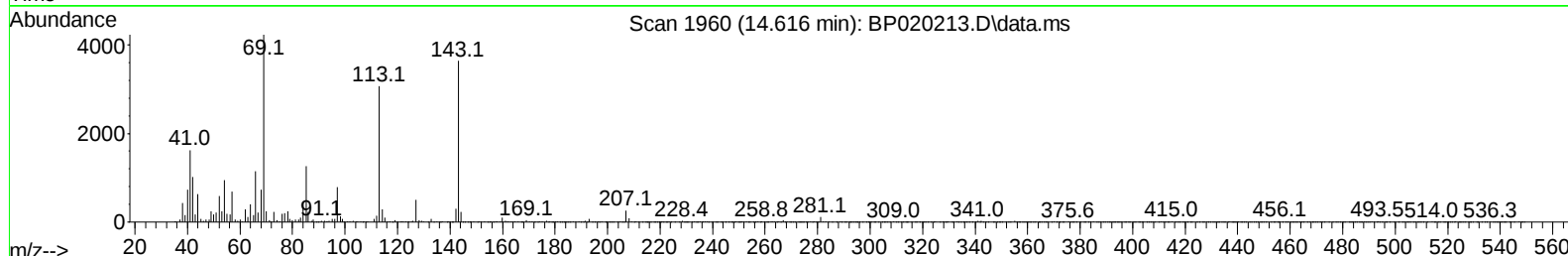
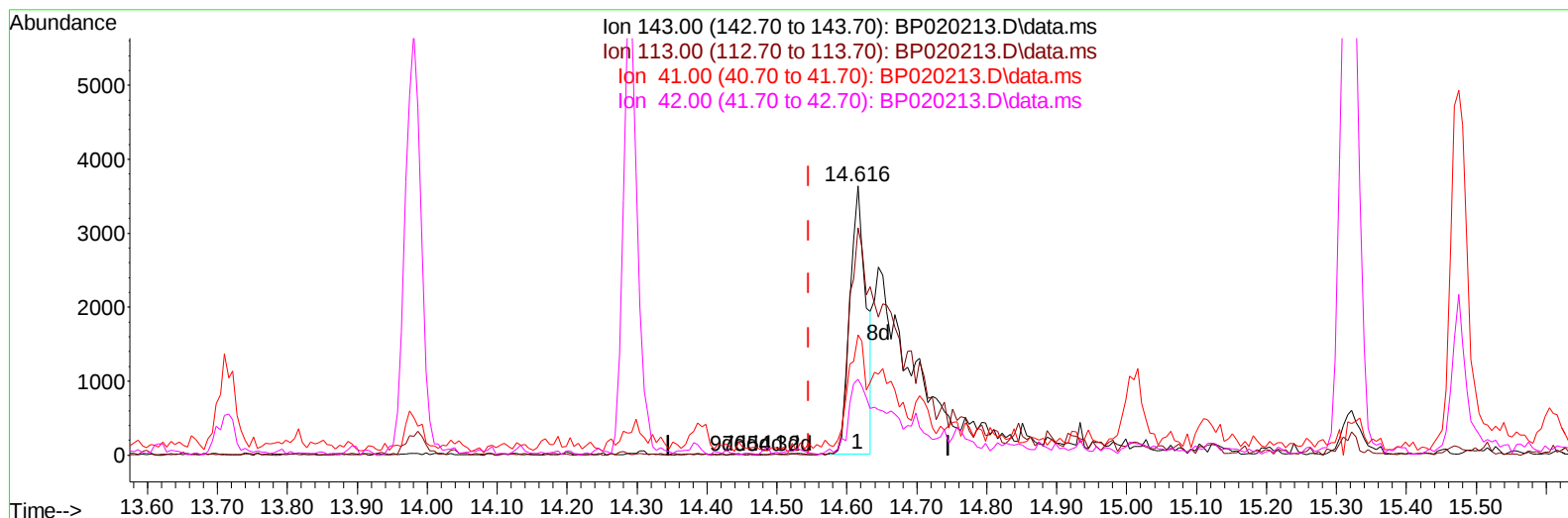
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050624\
 Data File : BP020213.D
 Acq On : 06 May 2024 19:33
 Operator : MA/JU
 Sample : P2373-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43Y2

Manual IntegrationsAPPROVED

Quant Time: May 06 22:45:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 06 12:26:12 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/07/2024
 Supervised By :mohammad ahmed 05/08/2024



TIC: BP020213.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.071) 2.40 ng/ul

response 5863

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	84.43
41.00	50.30	44.58
42.00	32.50	28.08

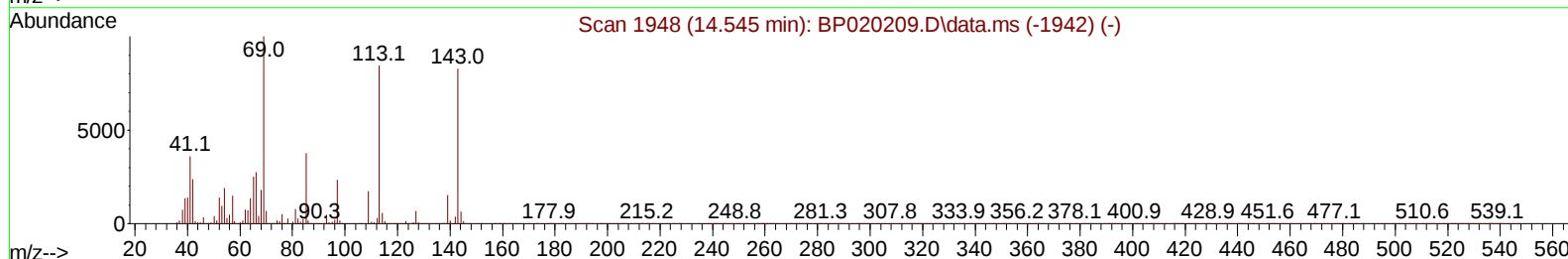
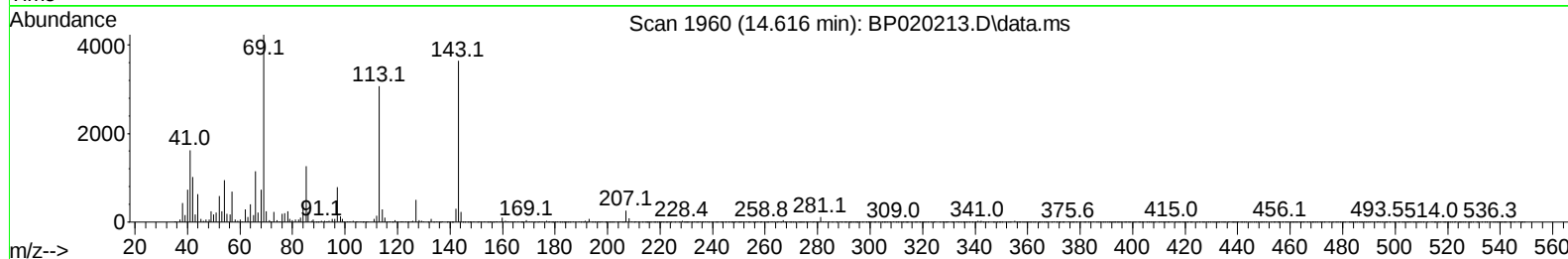
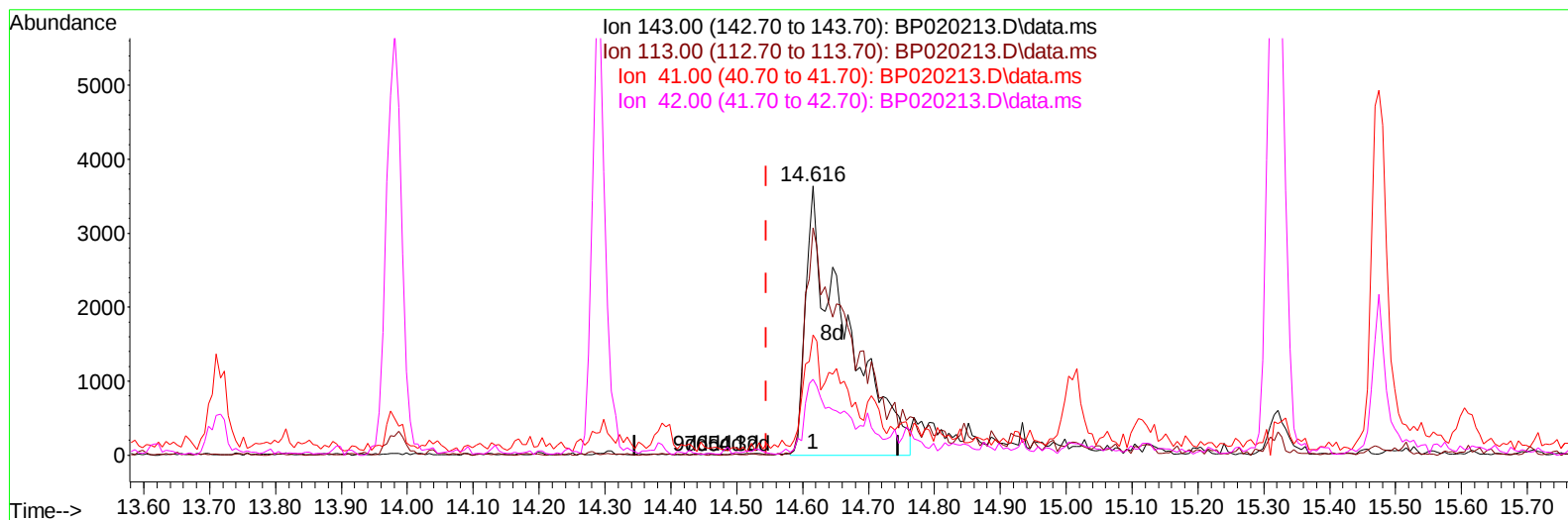
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Instrument :
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QLast Update : Mon May 06 12:26:12 2024
Response via : Initial Calibration



TIC: BP020213.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.071) 6.20 ng/ul m

response 15162

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	84.43
41.00	50.30	44.58
42.00	32.50	28.08

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050624\
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/07/2024
 Supervised By :mohammad ahmed 05/08/2024

Quant Time: May 07 03:29:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 06 12:26:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	65540	20.000	ng/ul	0.02
20) Naphthalene-d8	10.393	136	290351	20.000	ng/ul	0.02
38) Acenaphthene-d10	14.292	164	217767	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.128	188	530604	20.000	ng/ul	-0.01
79) Chrysene-d12	21.621	240	551493	20.000	ng/ul	-0.01
88) Perylene-d12	24.986	264	550023	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	6492	4.183	ng/uL	0.00
4) Pyridine-d5	3.634	84	23887	5.297	ng/ul	0.03
7) Phenol-d5	6.822	99	35003	6.219	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	6.975	67	107858	31.545	ng/ul	0.02
11) 2-Chlorophenol-d4	7.163	132	99708	24.780	ng/ul	0.02
15) 4-Methylphenol-d8	8.334	113	72725	15.982	ng/ul	0.02
21) Nitrobenzene-d5	8.781	128	69347	31.895	ng/ul	0.02
24) 2-Nitrophenol-d4	9.493	143	74944	30.100	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.034	165	124451	27.391	ng/ul	0.02
31) 4-Chloroaniline-d4	10.557	131	161458	25.390	ng/ul	0.02
46) Dimethylphthalate-d6	13.716	166	583384	36.685	ng/ul	0.01
49) Acenaphthylene-d8	13.981	160	579791	33.004	ng/ul	0.00
54) 4-Nitrophenol-d4	14.616	143	15162m	6.203	ng/ul	0.07
60) Fluorene-d10	15.316	176	483011	36.415	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	97273	28.890	ng/ul	0.00
73) Anthracene-d10	17.239	188	876603	38.189	ng/ul	0.00
81) Pyrene-d10	19.639	212	1144924	35.515	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.762	264	1015781	38.195	ng/ul	0.00

Target Compounds Qvalue

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

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