

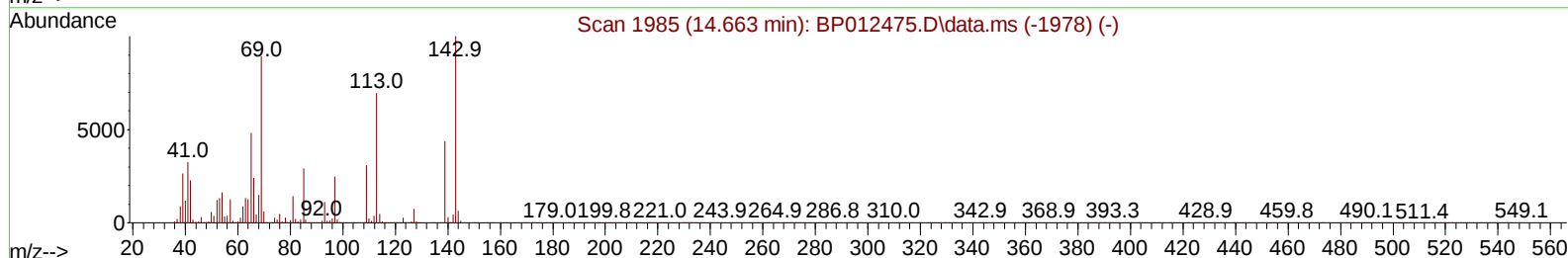
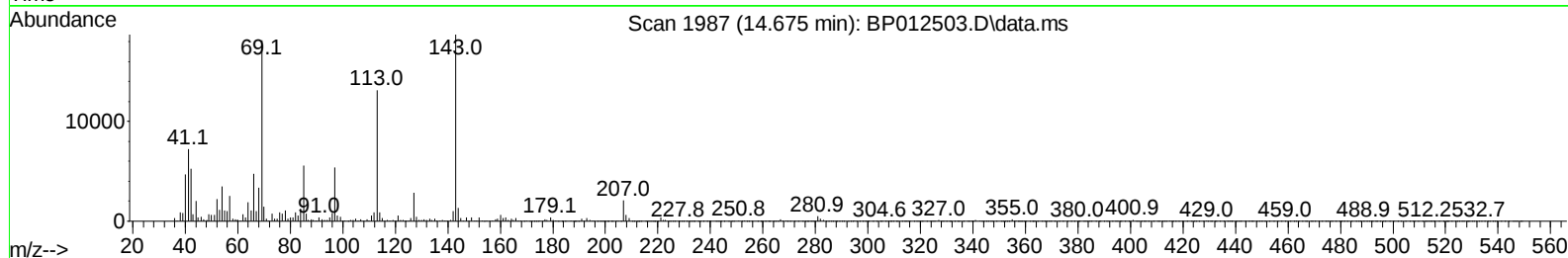
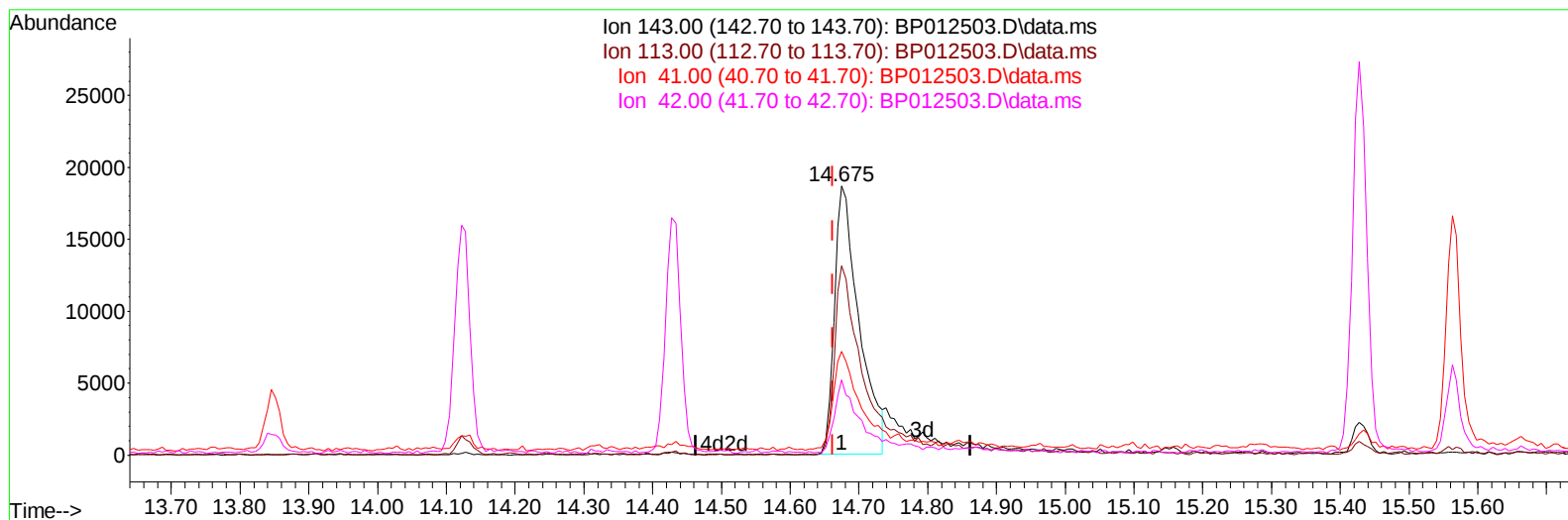
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012503.D
 Acq On : 03 Nov 2022 23:00
 Operator : CG/JU
 Sample : N5319-14
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAA8

Manual IntegrationsAPPROVED

Quant Time: Nov 04 01:04:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/05/2022
 Supervised By :mohammad ahmed 11/06/2022



TIC: BP012503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.012) 3.52 ng/u1

response 46387

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	70.28
41.00	34.50	38.59
42.00	23.60	27.98

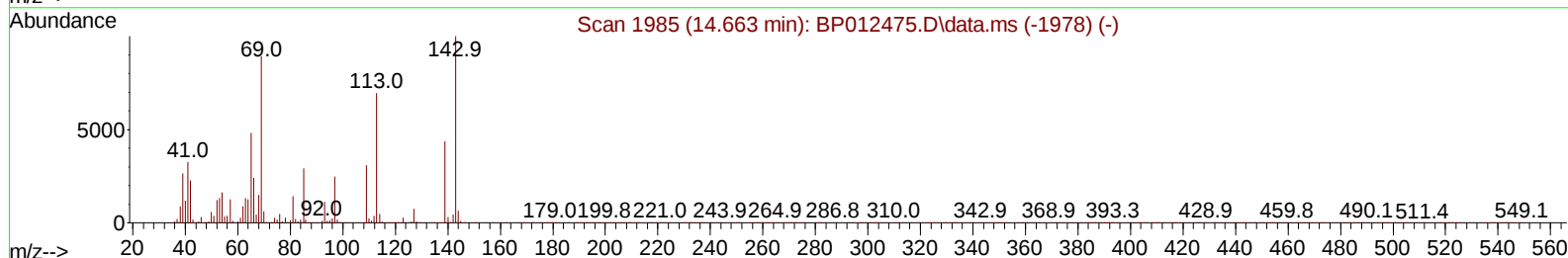
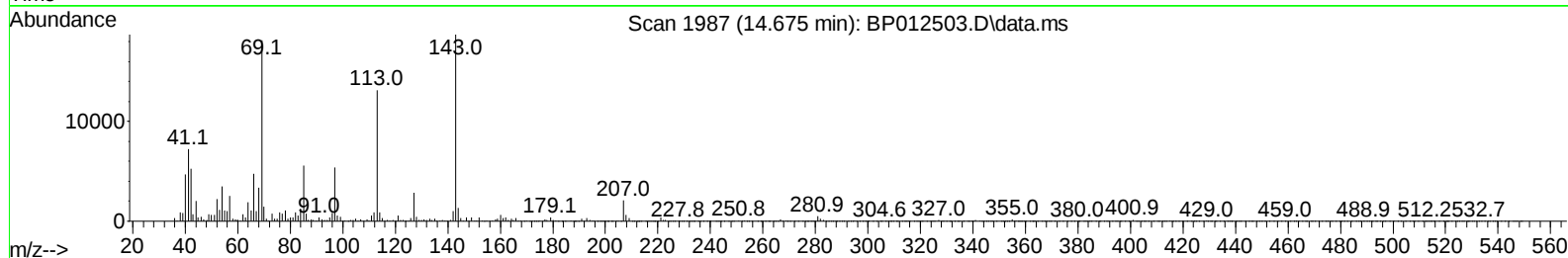
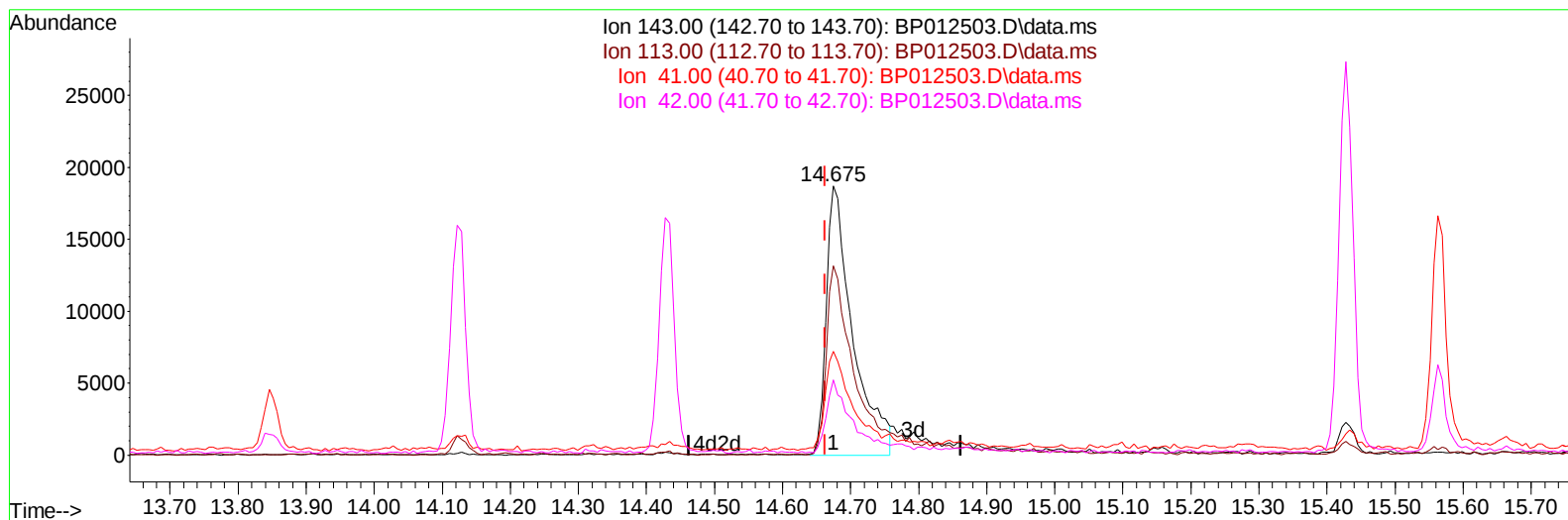
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 Sample : N5319-14
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Quant Time: Nov 04 01:04:05 2022
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TIC: BP012503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.012) 3.83 ng/ul m

response 50412

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	70.28
41.00	34.50	38.59
42.00	23.60	27.98

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 Operator : CG/JU
 Sample : N5319-14
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAA8

Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	312912	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1465644	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.434	164	1015832	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2304126	20.000	ng/ul	-0.01
79) Chrysene-d12	21.286	240	2136779	20.000	ng/ul	-0.01
88) Perylene-d12	23.669	264	1906658	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	28015	3.605	ng/uL	0.00
4) Pyridine-d5	3.658	84	192746	8.634	ng/ul	0.00
7) Phenol-d5	6.958	99	169499	6.039	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	495059	29.548	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.311	132	494118	24.109	ng/ul	-0.01
15) 4-Methylphenol-d8	8.499	113	345324	15.557	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	339693	35.741	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	364813	36.981	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	633357	29.625	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.734	131	813307	25.550	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	2288684	32.912	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	2510725	31.594	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.675	143	50412m	3.830	ng/ul	0.01
60) Fluorene-d10	15.428	176	1978101	33.224	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	324332	27.036	ng/ul	-0.01
73) Anthracene-d10	17.292	188	3325468	33.496	ng/ul	-0.01
81) Pyrene-d10	19.534	212	3830869	35.709	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.516	264	3196916	34.316	ng/ul	-0.01

Target Compounds Qvalue

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

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