

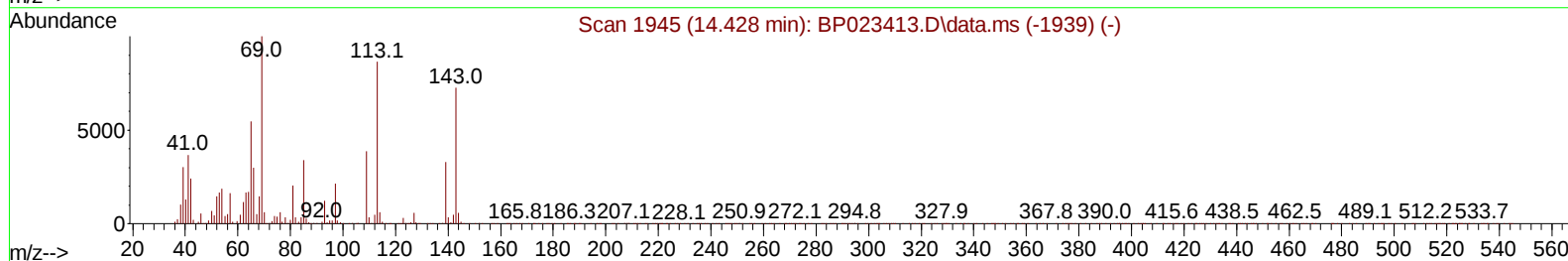
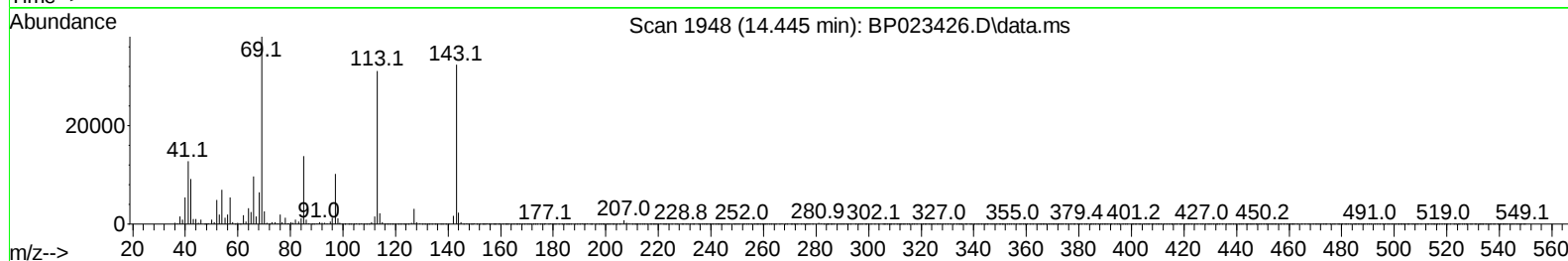
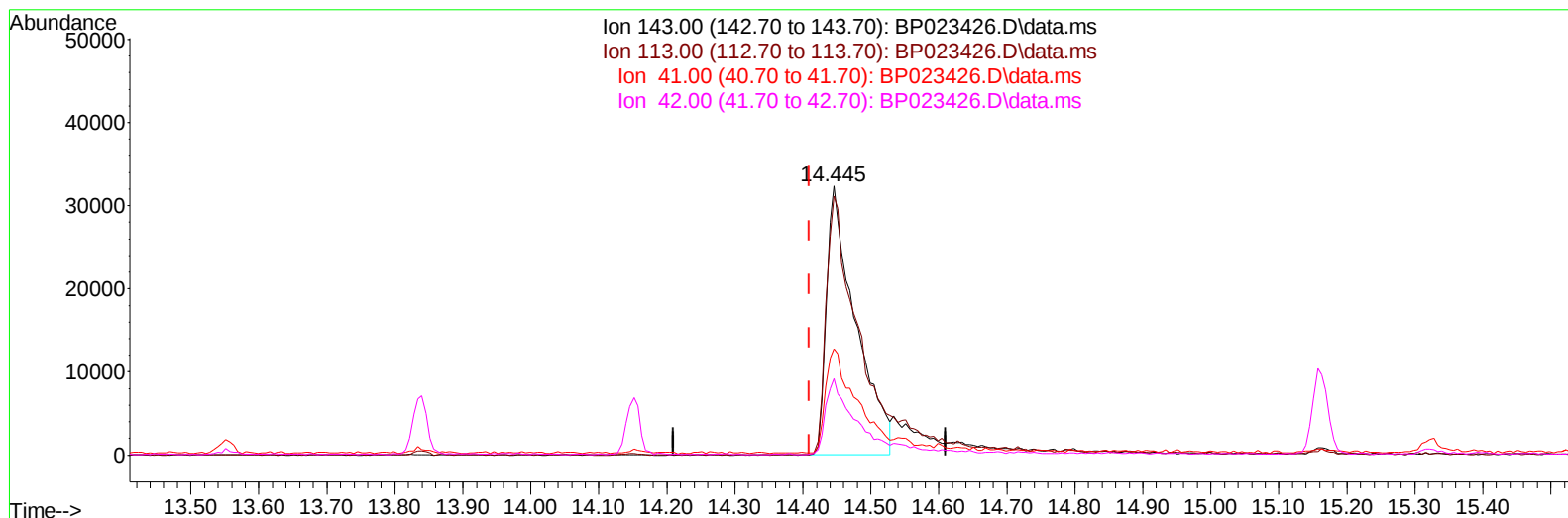
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121324\
Data File : BP023426.D
Acq On : 13 Dec 2024 18:52
Operator : RC/JU
Sample : PB165624BL
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK624

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/14/2024
Supervised By :mohammad ahmed 12/14/2024

Quant Time: Dec 13 22:49:22 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration



TIC: BP023426.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.445min (+ 0.035) 29.86 ng/u1

response 96179

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	96.36
41.00	45.60	39.40
42.00	31.70	28.38

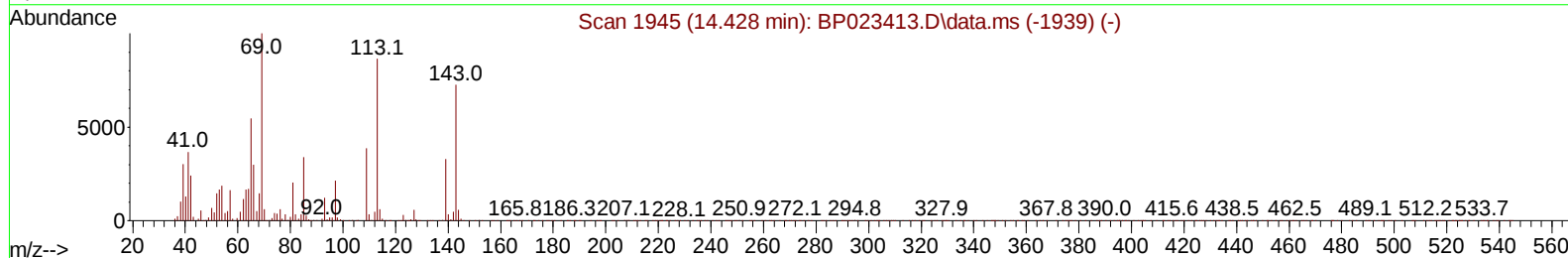
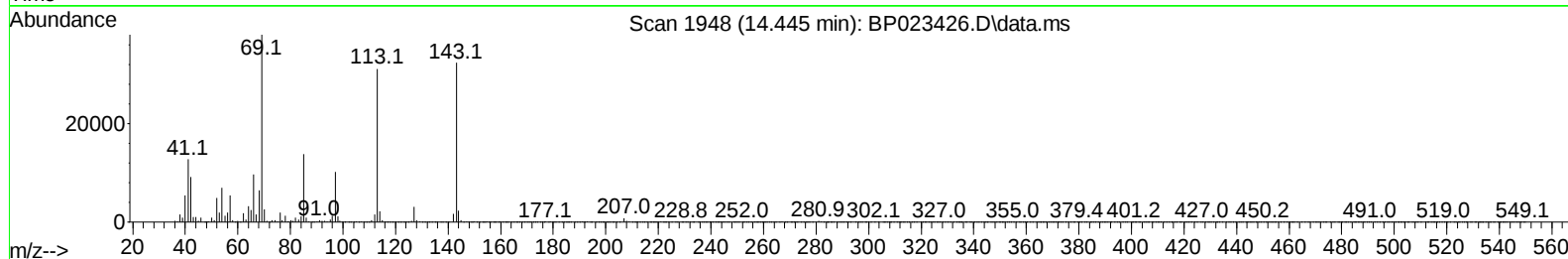
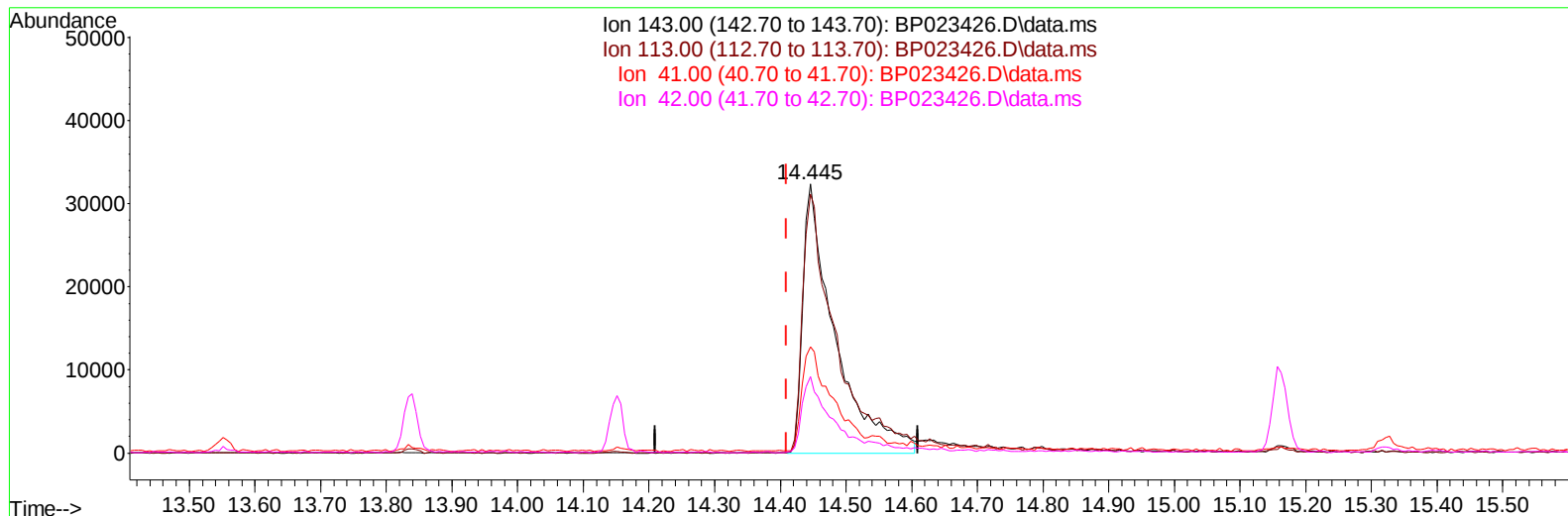
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TIC: BP023426.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.445min (+ 0.035) 33.91 ng/ul m

response 109210

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	96.36
41.00	45.60	39.40
42.00	31.70	28.38

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/14/2024
 Supervised By :mohammad ahmed 12/14/2024

Quant Time: Dec 13 22:51:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.534	152	109729	20.000	ng/ul	0.00
20) Naphthalene-d8	10.281	136	423454	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.151	164	338585	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.951	188	783437	20.000	ng/ul	0.00
79) Chrysene-d12	21.363	240	820606	20.000	ng/ul	-0.02
88) Perylene-d12	24.527	264	879488	20.000	ng/ul	-0.05
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	14570	4.889	ng/uL	0.00
4) Pyridine-d5	3.511	84	215638	28.136	ng/ul	0.00
7) Phenol-d5	6.752	99	270410	31.408	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.887	67	171829	31.633	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	198593	31.253	ng/ul	0.00
15) 4-Methylphenol-d8	8.252	113	217499	31.301	ng/ul	0.00
21) Nitrobenzene-d5	8.681	128	98986	31.304	ng/ul	0.00
24) 2-Nitrophenol-d4	9.387	143	111090	30.483	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.934	165	234489	30.486	ng/ul	0.00
31) 4-Chloroaniline-d4	10.434	131	260705	30.256	ng/ul	0.00
46) Dimethylphthalate-d6	13.551	166	813737	30.959	ng/ul	0.00
49) Acenaphthylene-d8	13.840	160	864227	31.172	ng/ul	0.00
54) 4-Nitrophenol-d4	14.445	143	109210m	33.906	ng/ul	0.04
60) Fluorene-d10	15.163	176	720046	31.145	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.322	200	43900	9.272	ng/ul	0.01
73) Anthracene-d10	17.051	188	1255774	35.122	ng/ul	0.00
81) Pyrene-d10	19.439	212	1528385	35.649	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.304	264	1490617	31.661	ng/ul	-0.04

Target Compounds Qvalue

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

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