

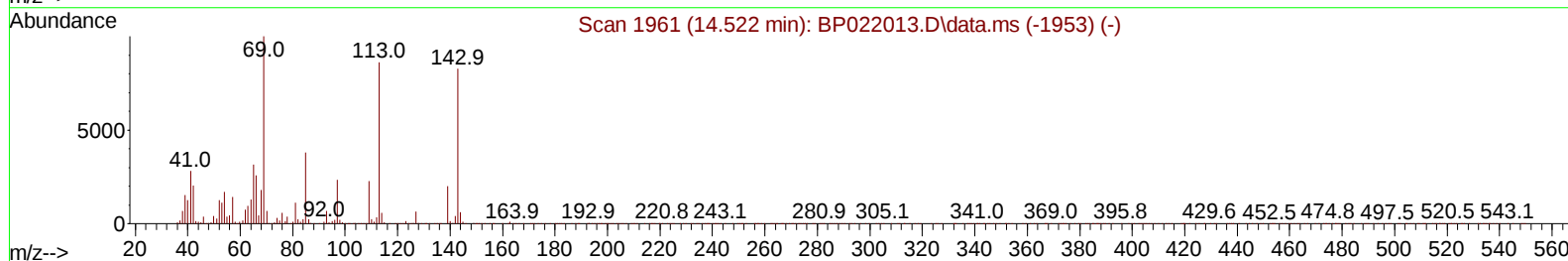
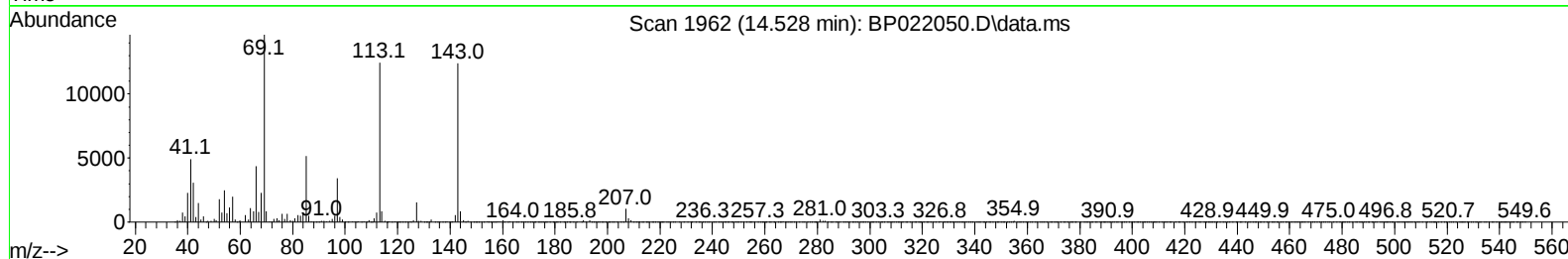
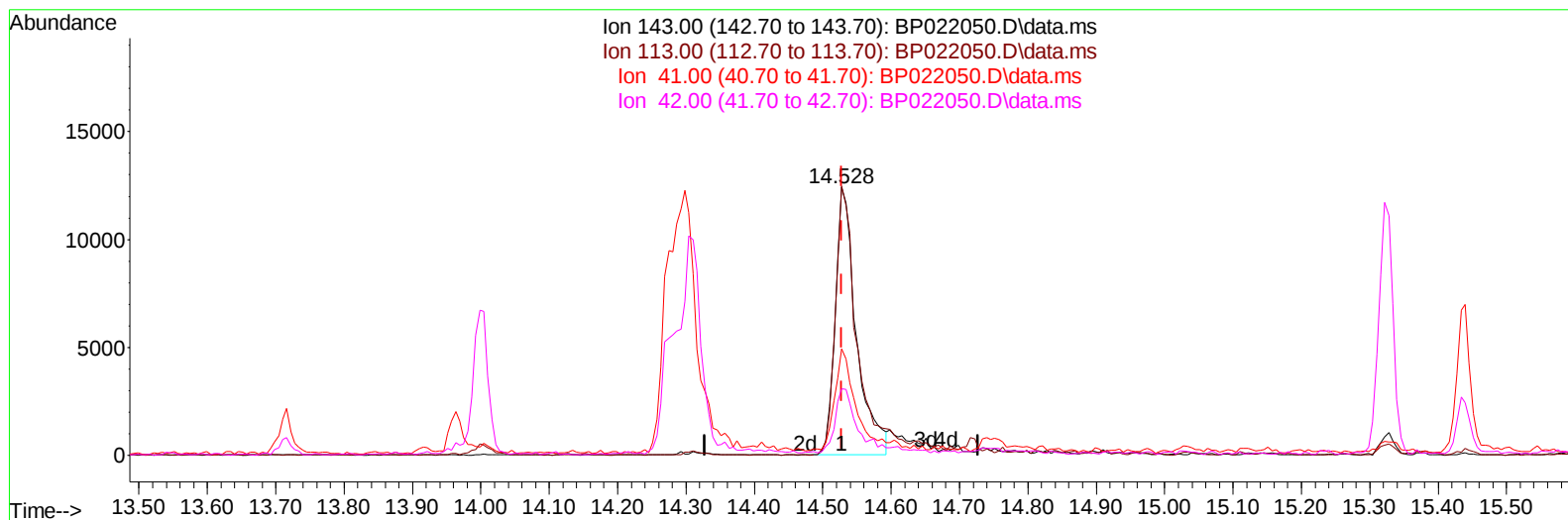
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022050.D
Acq On : 26 Sep 2024 06:21
Operator : RC/JU
Sample : P4107-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B23

Manual IntegrationsAPPROVED

Quant Time: Sep 26 06:57:45 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2024
Supervised By :mohammad ahmed 09/28/2024



TIC: BP022050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.000) 6.29 ng/u1

response 26652

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	100.49
41.00	35.80	39.75
42.00	26.20	24.87

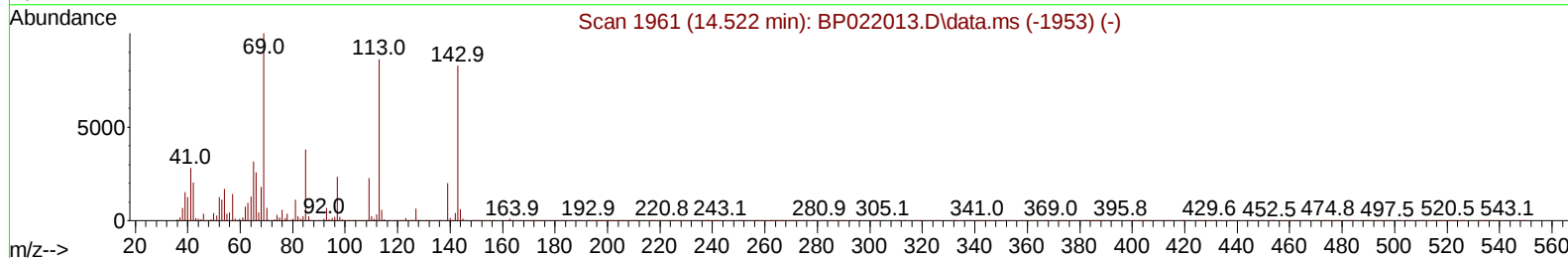
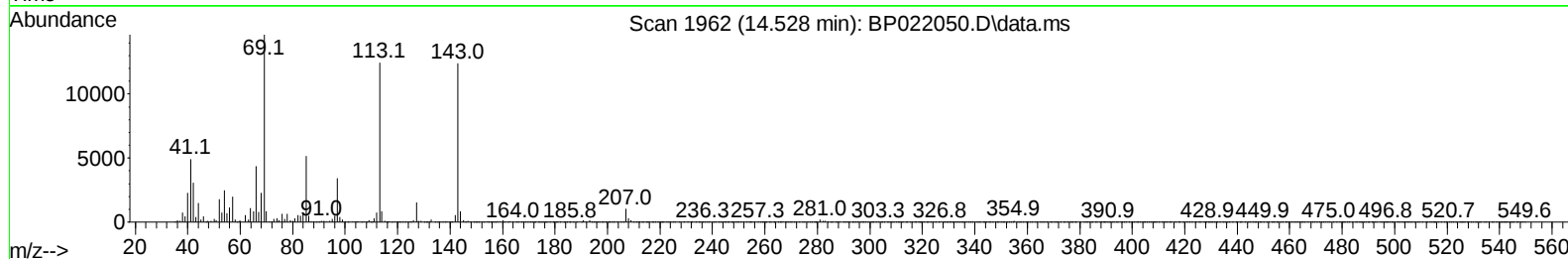
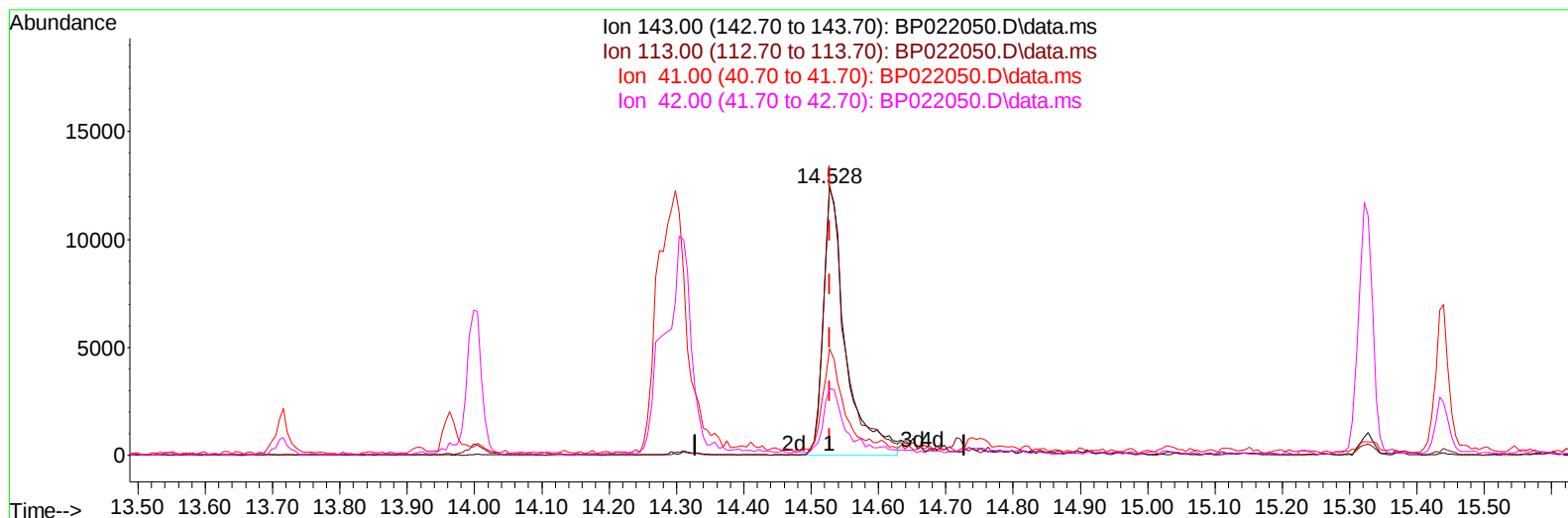
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TIC: BP022050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.000) 6.73 ng/ul m

response 28523

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	100.49
41.00	35.80	39.75
42.00	26.20	24.87

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/27/2024
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Quant Time: Sep 26 06:59:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	109873	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	422856	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	320675	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	791028	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	904530	20.000	ng/ul	0.00
88) Perylene-d12	24.827	264	995308	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	12892	4.596	ng/uL	0.00
4) Pyridine-d5	3.640	84	84035	10.481	ng/ul	0.00
7) Phenol-d5	6.893	99	83145	9.251	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	174327	31.499	ng/ul	0.00
11) 2-Chlorophenol-d4	7.252	132	197050	30.303	ng/ul	0.00
15) 4-Methylphenol-d8	8.405	113	163544	22.964	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	106086	33.519	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	137733	38.381	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	282510	36.251	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	232722	25.028	ng/ul	0.00
46) Dimethylphthalate-d6	13.716	166	925454	37.334	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	902456	34.386	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	28523m	6.727	ng/ul	0.00
60) Fluorene-d10	15.328	176	794633	36.240	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	166330	33.868	ng/ul	0.00
73) Anthracene-d10	17.210	188	1458869	39.505	ng/ul	0.00
81) Pyrene-d10	19.586	212	1946972	41.547	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.604	264	2108613	40.279	ng/ul	0.00

Target Compounds Qvalue

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

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