

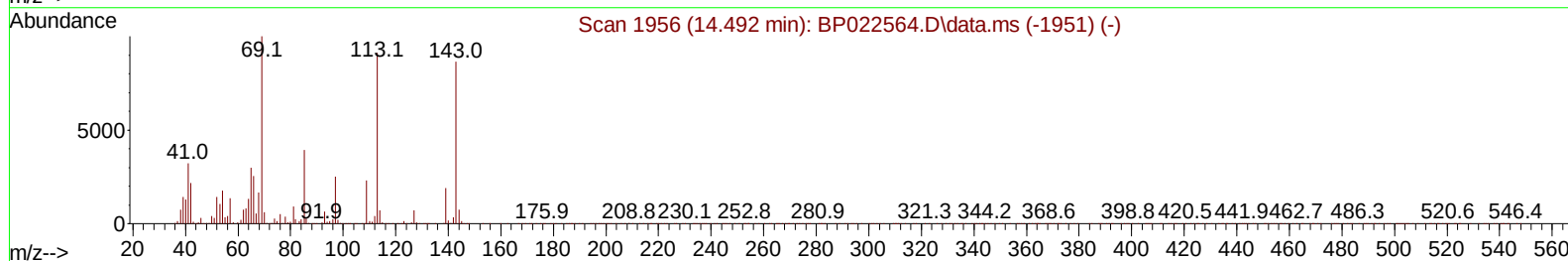
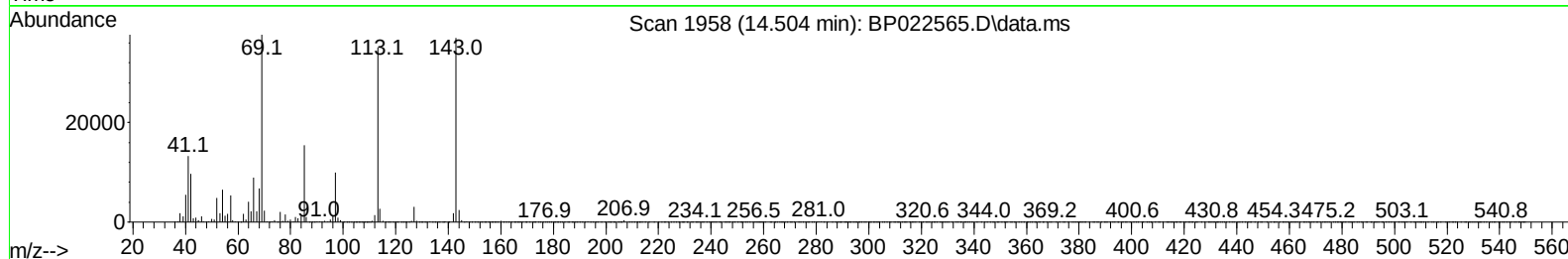
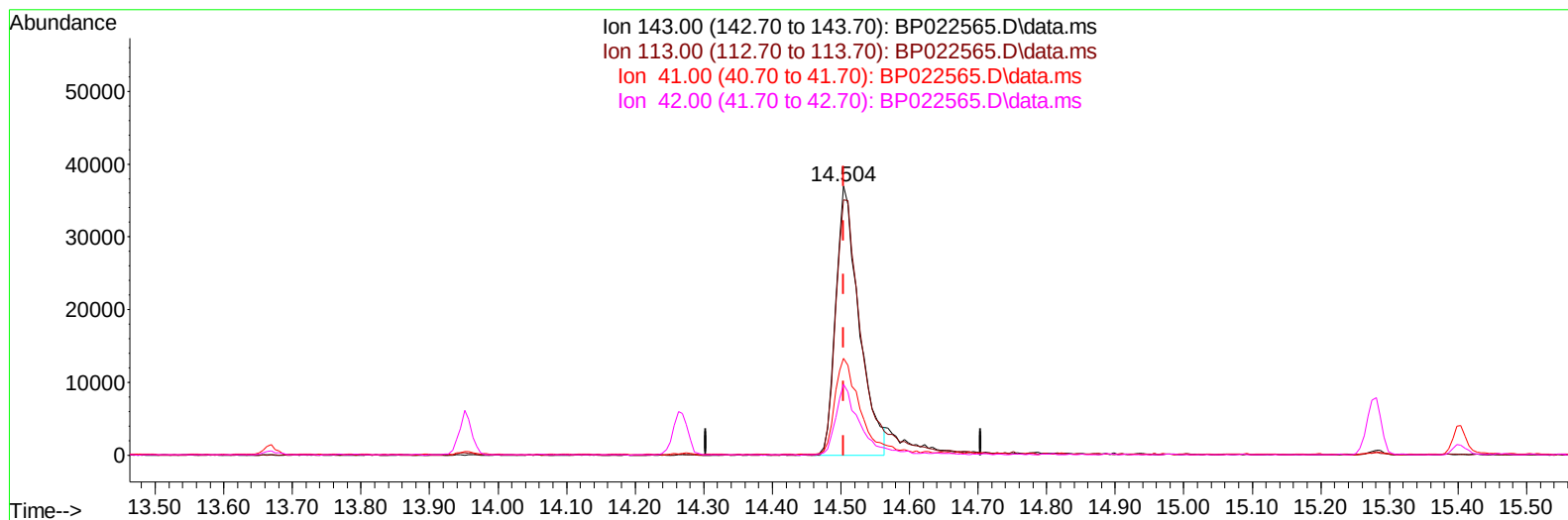
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022565.D
Acq On : 23 Oct 2024 16:11
Operator : RC/JU
Sample : PB164327BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK327

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/24/2024
Supervised By :mohammad ahmed 10/25/2024

Quant Time: Oct 23 17:10:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 22 05:59:59 2024
Response via : Initial Calibration



TIC: BP022565.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 19.94 ng/ul

response 86864

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	94.98
41.00	35.90	35.87
42.00	23.70	26.35

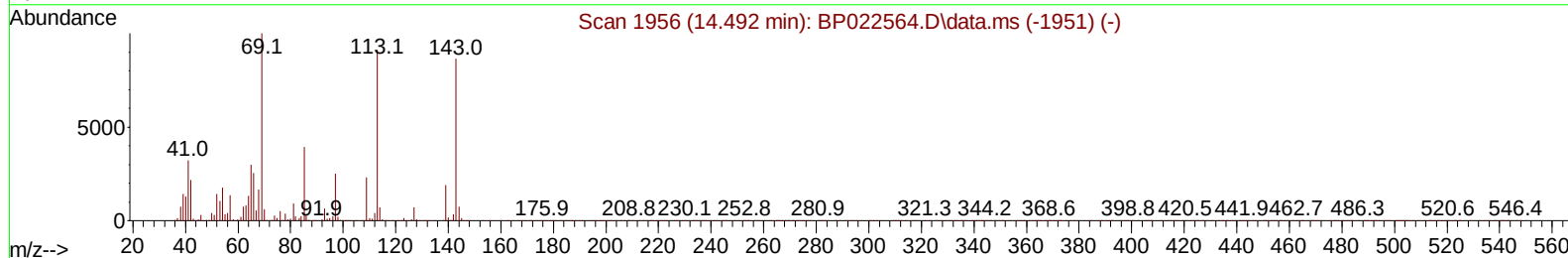
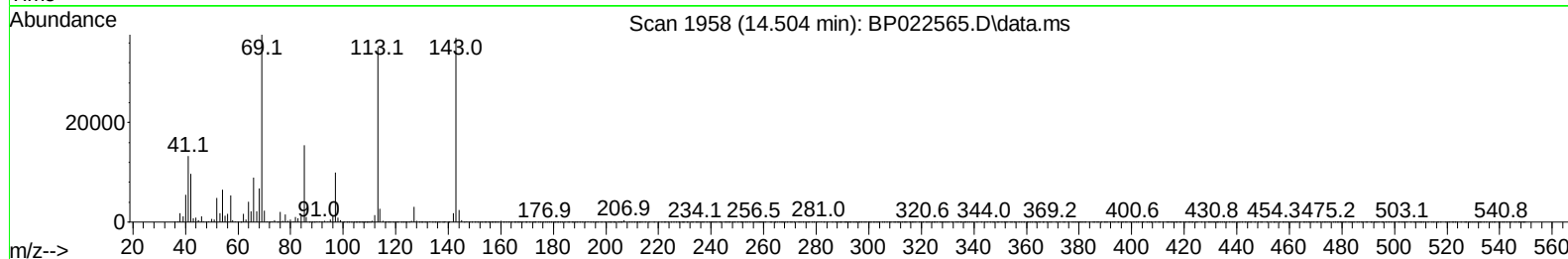
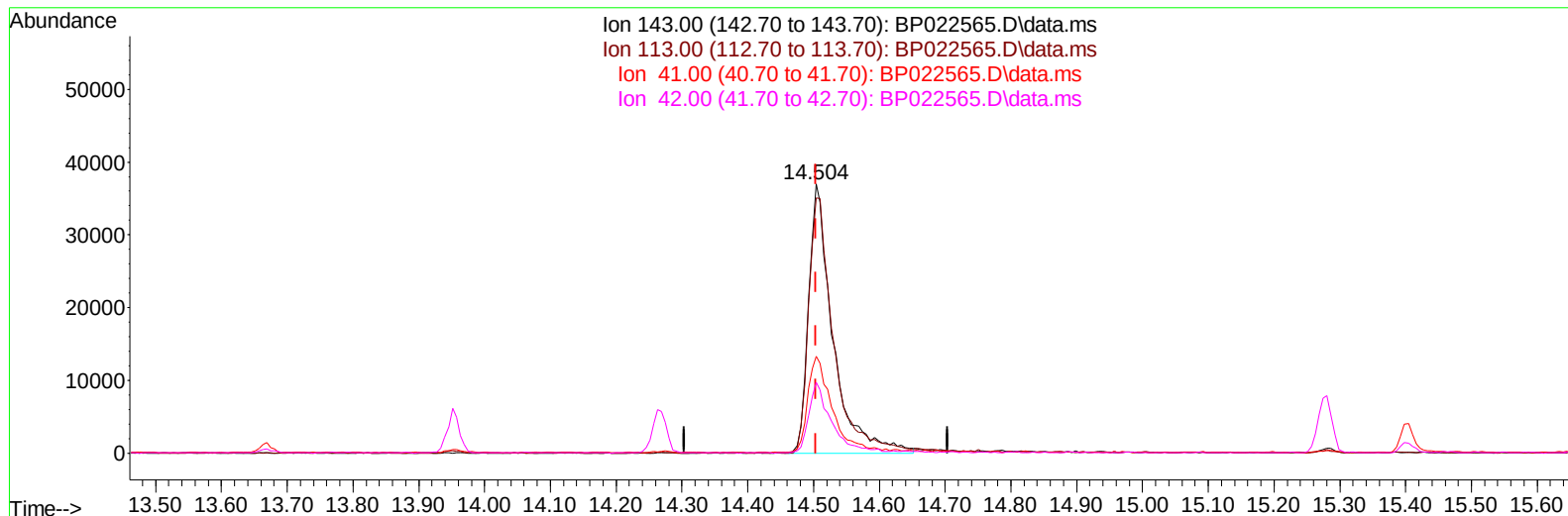
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TIC: BP022565.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 21.90 ng/ul m

response 95405

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	94.98
41.00	35.90	35.87
42.00	23.70	26.35

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Quant Time: Oct 23 17:26:49 2024
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 QLast Update : Tue Oct 22 05:59:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	112961	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	429244	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	326881	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	804132	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	807794	20.000	ng/ul	0.00
88) Perylene-d12	24.774	264	875193	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	12776	4.395	ng/uL	0.00
4) Pyridine-d5	3.605	84	179180	22.453	ng/ul	0.00
7) Phenol-d5	6.852	99	222776	23.429	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.999	67	132685	23.742	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	170057	25.204	ng/ul	0.00
15) 4-Methylphenol-d8	8.357	113	181453	23.912	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	83187	25.205	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	99979	26.165	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	197500	24.319	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.551	131	230924	24.064	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	656376	25.273	ng/ul	-0.01
49) Acenaphthylene-d8	13.957	160	695009	25.195	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	95405m	21.897	ng/ul	0.00
60) Fluorene-d10	15.281	176	567273	25.182	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	83478	15.784	ng/ul	-0.01
73) Anthracene-d10	17.169	188	964467	25.496	ng/ul	0.00
81) Pyrene-d10	19.557	212	1190010	27.858	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.557	264	1229487	26.305	ng/ul	0.03

Target Compounds Qvalue

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

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