

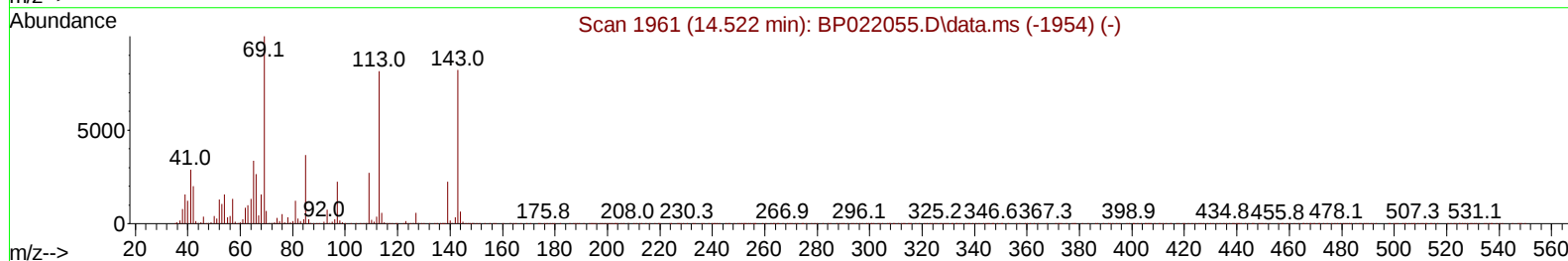
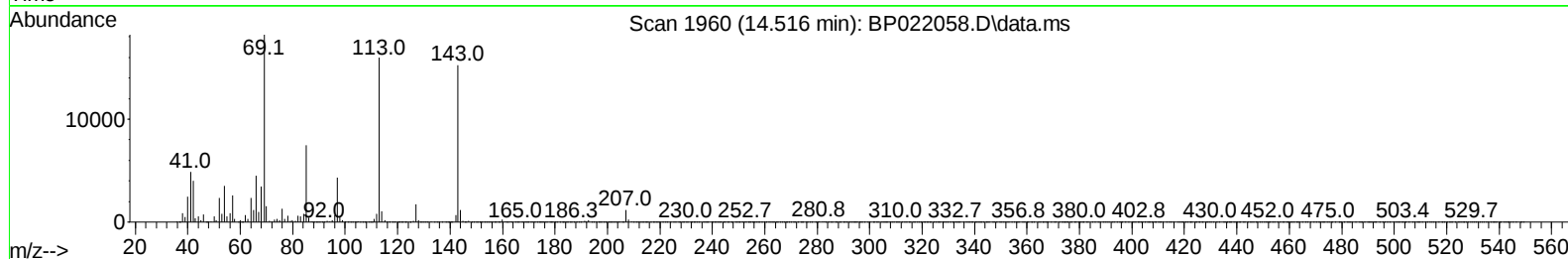
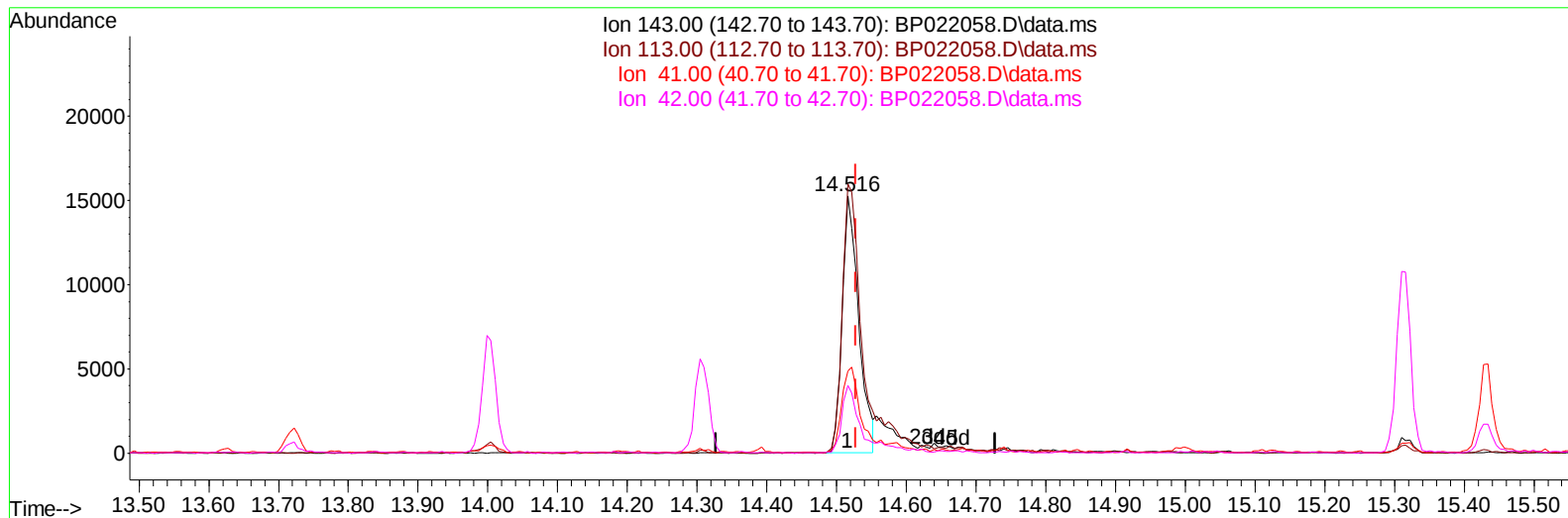
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022058.D
Acq On : 26 Sep 2024 11:47
Operator : RC/JU
Sample : P4084-06
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYM3

Manual IntegrationsAPPROVED

Quant Time: Sep 27 01:16:30 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: BP022058.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (-0.012) 8.00 ng/u1

response 24951

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	104.60
41.00	35.80	31.84
42.00	26.20	26.39

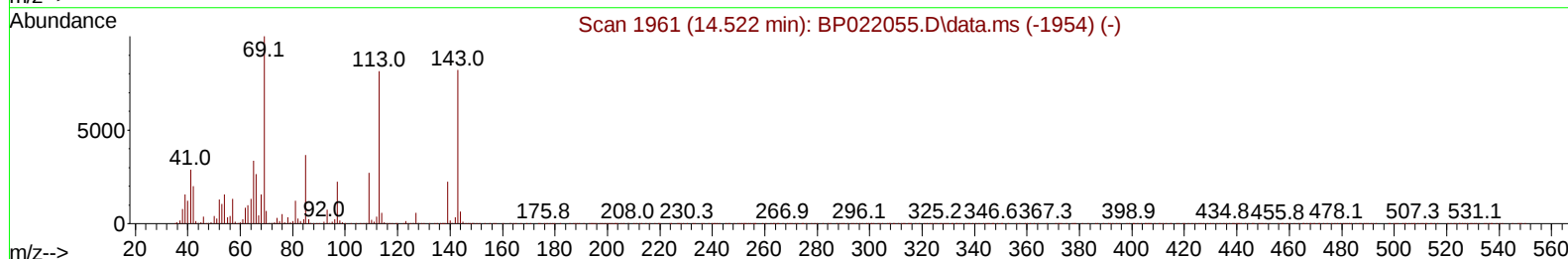
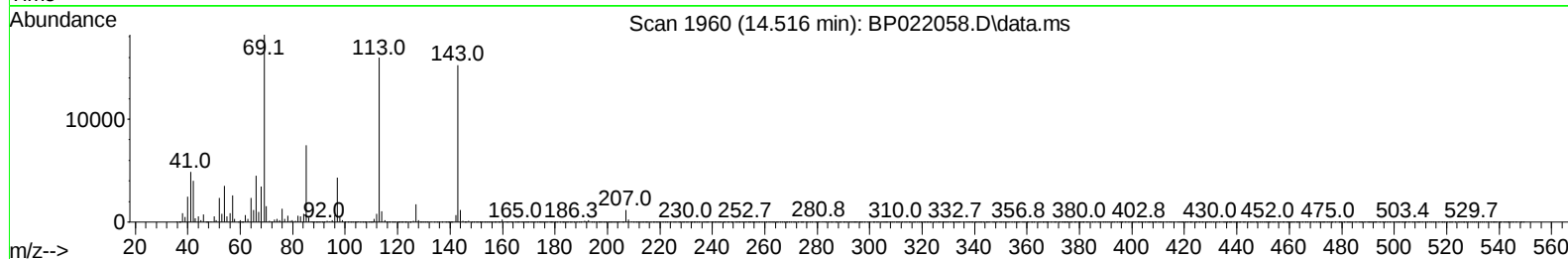
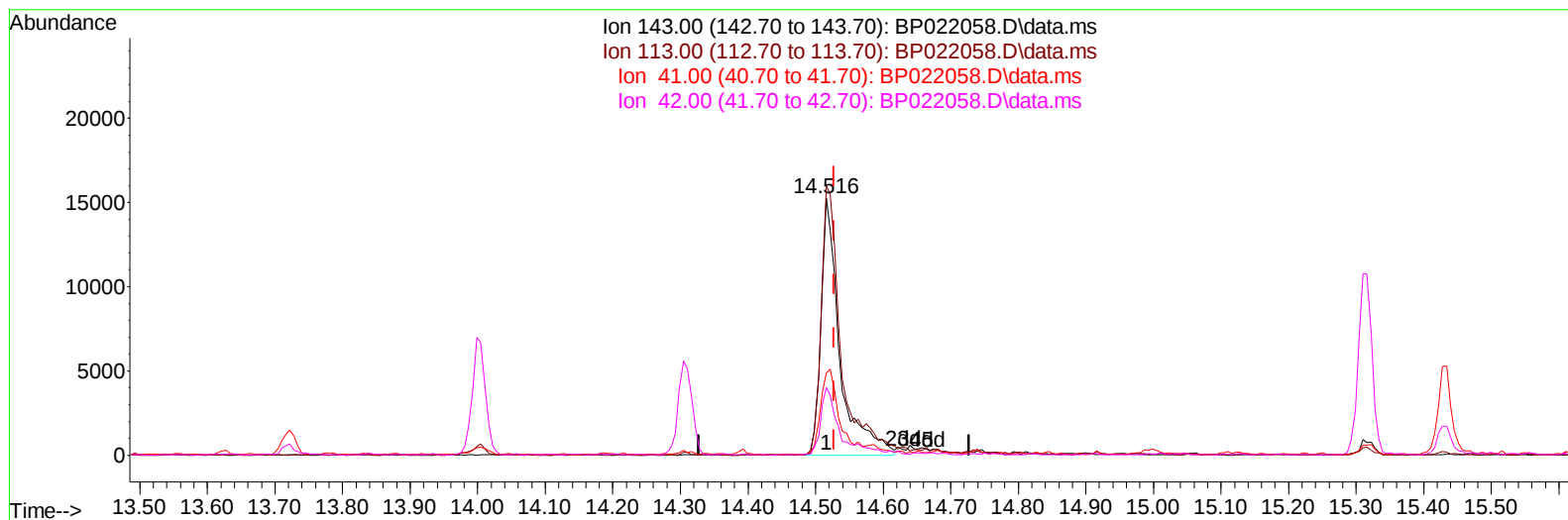
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TIC: BP022058.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (-0.012) 9.50 ng/u1 m

response 29645

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	104.60
41.00	35.80	31.84
42.00	26.20	26.39

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 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
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 DCYM3

Manual IntegrationsAPPROVED

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

Quant Time: Sep 27 01:23:43 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	7.705	152	84394	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	313933	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.304	164	235972	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.098	188	580735	20.000	ng/ul	-0.02
79) Chrysene-d12	21.528	240	732828	20.000	ng/ul	-0.01
88) Perylene-d12	24.804	264	885375	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	12792	5.937	ng/uL	0.00
4) Pyridine-d5	3.640	84	39383	6.395	ng/ul	0.00
7) Phenol-d5	6.887	99	71879	10.412	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	139885	32.907	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	160150	32.063	ng/ul	0.00
15) 4-Methylphenol-d8	8.405	113	122109	22.322	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	83829	35.677	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	100997	37.909	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.093	165	198532	34.314	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	124805	18.079	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	713458	39.113	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	681058	35.265	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	29645m	9.501	ng/ul	-0.01
60) Fluorene-d10	15.316	176	606857	37.611	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.434	200	119947	33.268	ng/ul	-0.01
73) Anthracene-d10	17.198	188	1081778	39.902	ng/ul	-0.02
81) Pyrene-d10	19.575	212	1486900	39.164	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.586	264	1844899	39.618	ng/ul	-0.02

Target Compounds Qvalue

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

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