

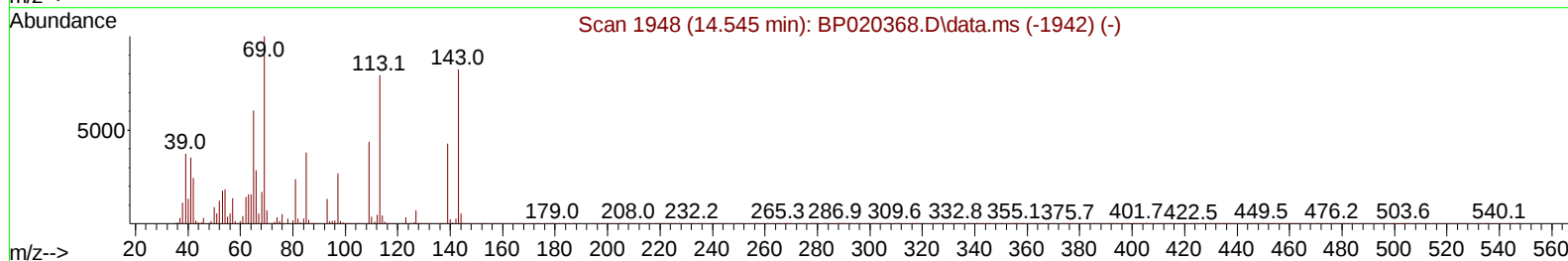
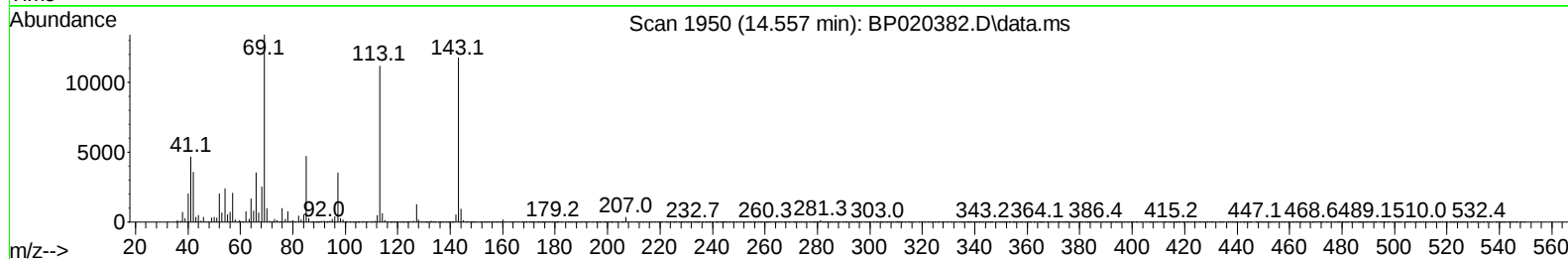
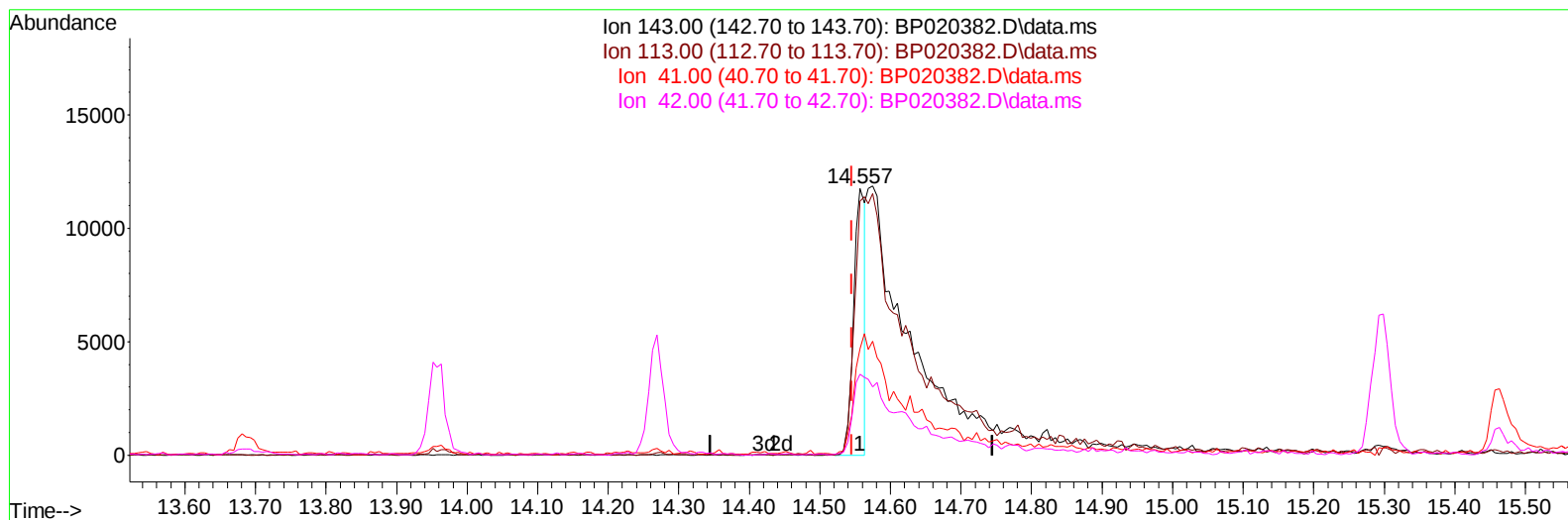
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051424\
 Data File : BP020382.D
 Acq On : 15 May 2024 01:41
 Operator : MA/JU
 Sample : PB160788BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK788

Manual IntegrationsAPPROVED

Quant Time: May 15 02:12:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 14 23:58:44 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024



TIC: BP020382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.012) 5.13 ng/u1

response 13518

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	95.17
41.00	50.30	39.85#
42.00	32.50	30.44

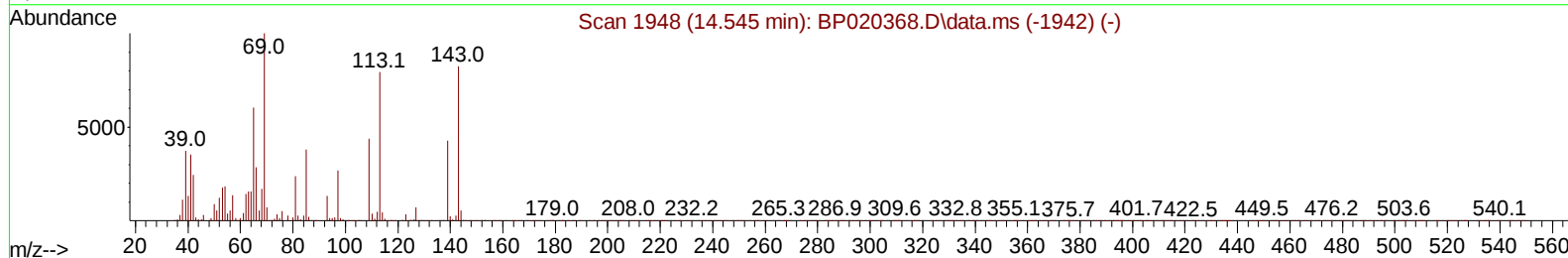
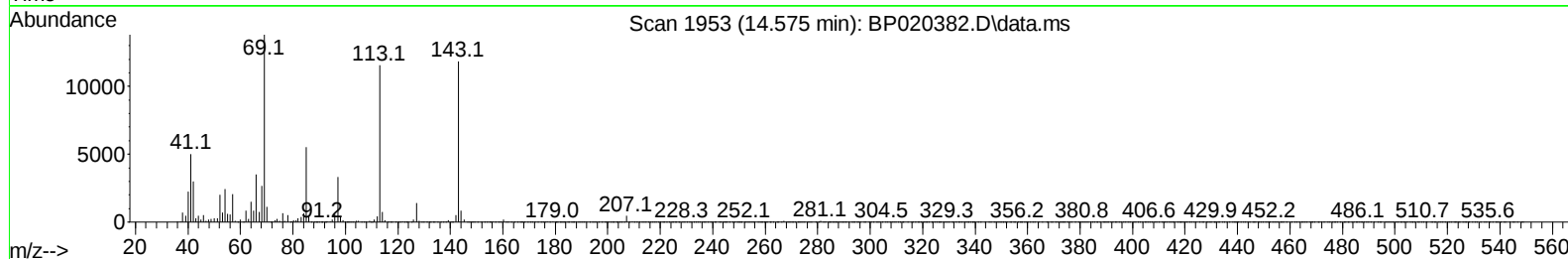
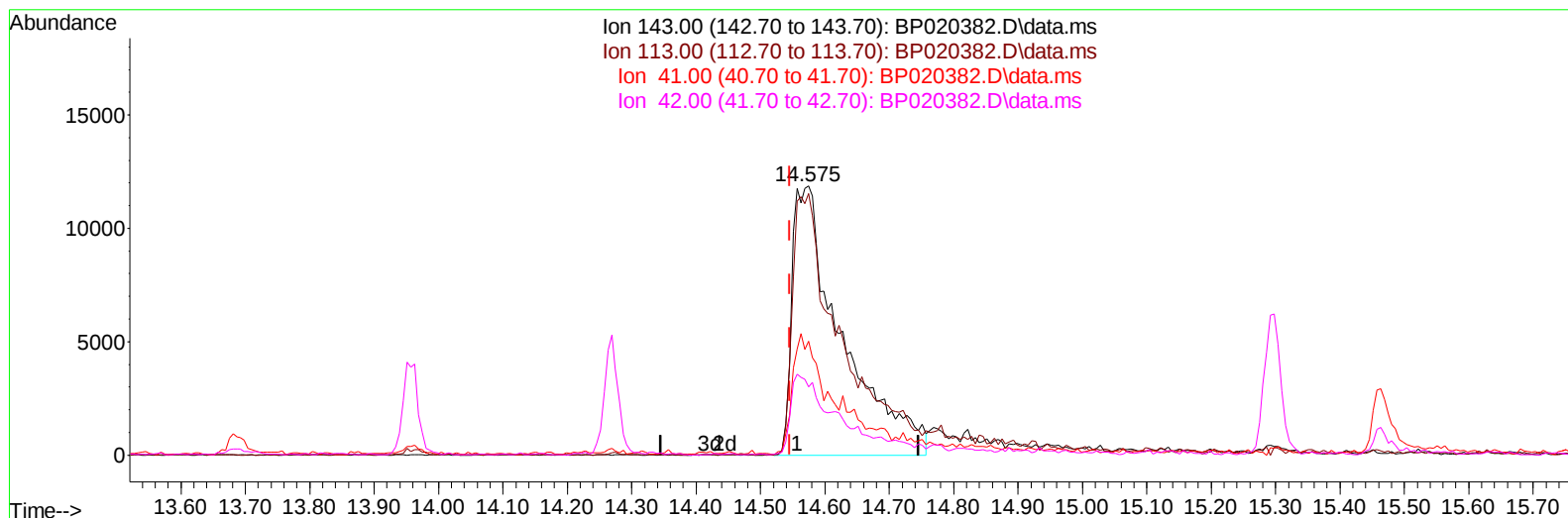
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TIC: BP020382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.029) 24.02 ng/ul m

response 63242

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.36
41.00	50.30	42.40
42.00	32.50	25.43#

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 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024

Quant Time: May 15 02:13:25 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 14 23:58:44 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	84456	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	341781	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	234566	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	545109	20.000	ng/ul	0.00
79) Chrysene-d12	21.604	240	601775	20.000	ng/ul	0.00
88) Perylene-d12	24.951	264	685995	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	11652	5.827	ng/uL	0.00
4) Pyridine-d5	3.599	84	145180	24.985	ng/ul	0.00
7) Phenol-d5	6.787	99	180554	24.894	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	110129	24.995	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	145919	28.143	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	144661	24.670	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	70089	27.385	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	78898	26.920	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	123079	23.012	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	180687	24.138	ng/ul	0.01
46) Dimethylphthalate-d6	13.687	166	489703	28.588	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	513882	27.158	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	63242m	24.022	ng/ul	0.03
60) Fluorene-d10	15.292	176	407857	28.547	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	77127	22.297	ng/ul	0.00
73) Anthracene-d10	17.210	188	669718	28.400	ng/ul	0.00
81) Pyrene-d10	19.622	212	841576	23.924	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.727	264	939958	28.339	ng/ul	0.01

Target Compounds Qvalue

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

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