

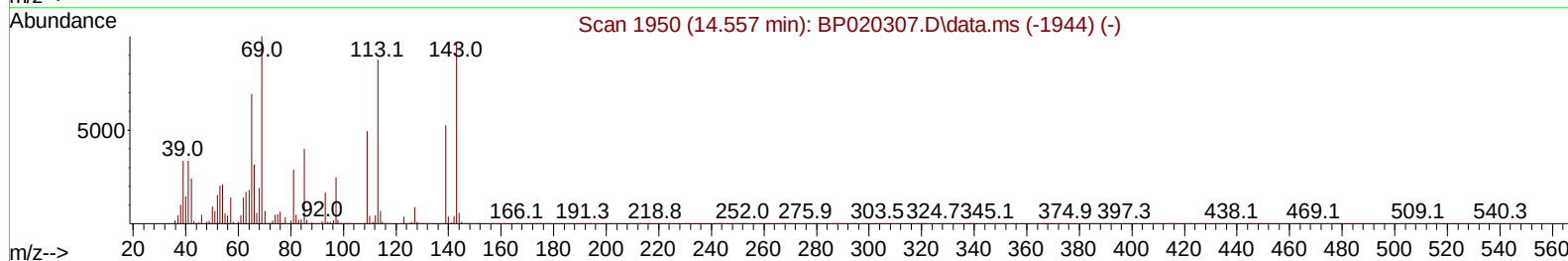
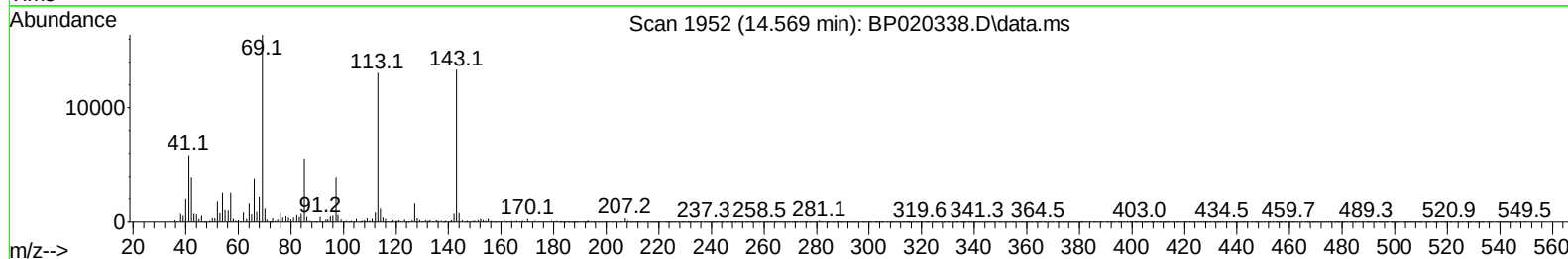
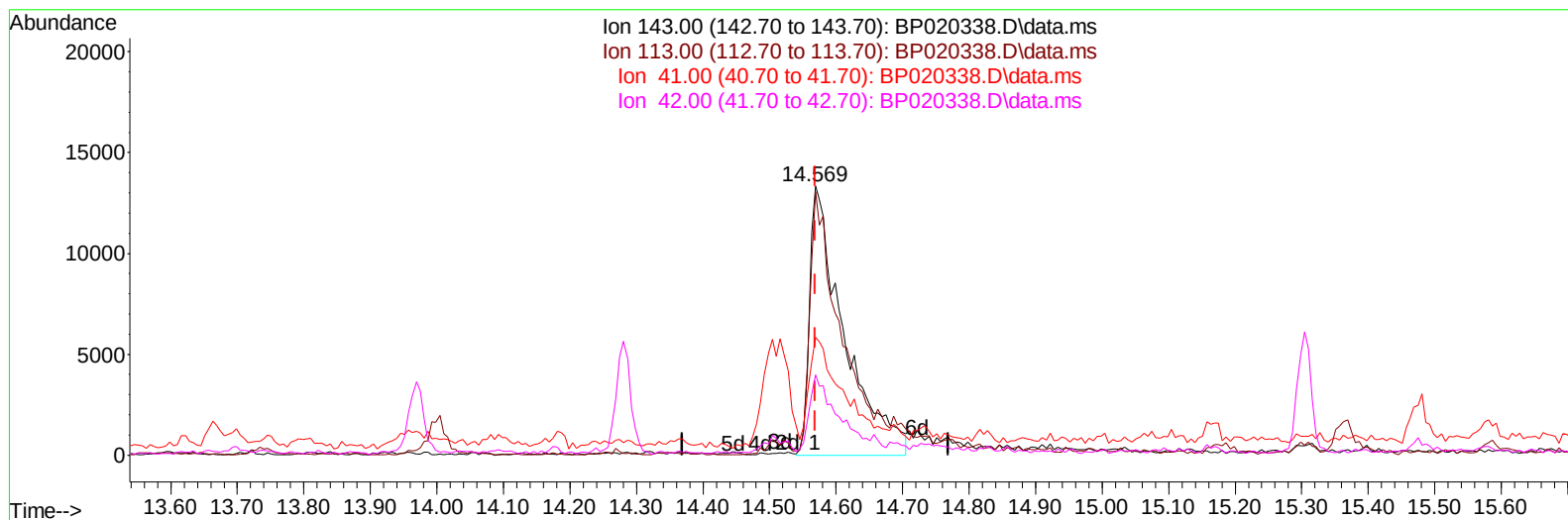
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020338.D
Acq On : 12 May 2024 03:23
Operator : MA/JU
Sample : P2371-05
Misc :
ALS Vial : 59 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R7

Manual IntegrationsAPPROVED

Quant Time: May 13 00:27:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon May 13 00:21:03 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/13/2024
Supervised By :mohammad ahmed 05/14/2024



TIC: BP020338.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.000) 15.76 ng/ul m

response 47626

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.89
41.00	50.30	43.77
42.00	32.50	29.76

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Reviewed By :Jagrut Upadhyay 05/13/2024
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Quant Time: May 13 01:10:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 13 00:21:03 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	95657	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	399746	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164	269259	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.128	188	579671	20.000	ng/ul	0.00
79) Chrysene-d12	21.616	240	581448	20.000	ng/ul	0.00
88) Perylene-d12	24.974	264	659467	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	6581	2.906	ng/uL	0.00
4) Pyridine-d5	3.628	84	52537	7.983	ng/ul	0.01
7) Phenol-d5	6.805	99	136484	16.614	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67	80833	16.198	ng/ul	0.00
11) 2-Chlorophenol-d4	7.152	132	107465	18.299	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	96234	14.490	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	54109	18.076	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	62439	18.215	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	112757	18.025	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	76508	8.739	ng/ul	0.01
46) Dimethylphthalate-d6	13.698	166	389957	19.832	ng/ul	-0.01
49) Acenaphthylene-d8	13.969	160	423648	19.504	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143	47626m	15.760	ng/ul	0.00
60) Fluorene-d10	15.304	176	338931	20.666	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	48076	13.070	ng/ul	0.00
73) Anthracene-d10	17.228	188	551875	22.007	ng/ul	0.00
81) Pyrene-d10	19.639	212	671697	19.762	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.751	264	684269	21.460	ng/ul	0.01

Target Compounds Qvalue

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

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