

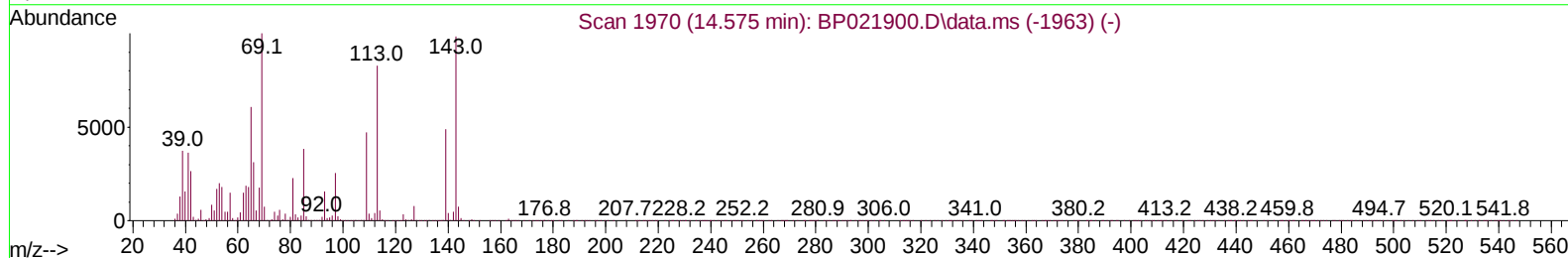
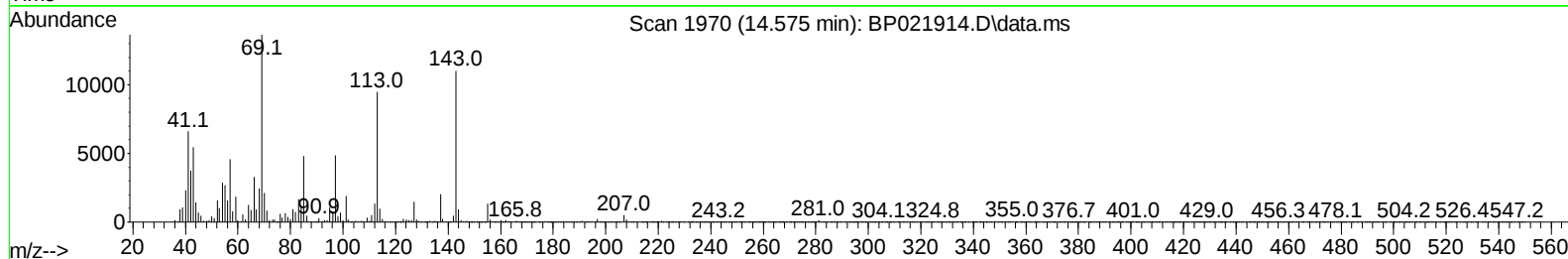
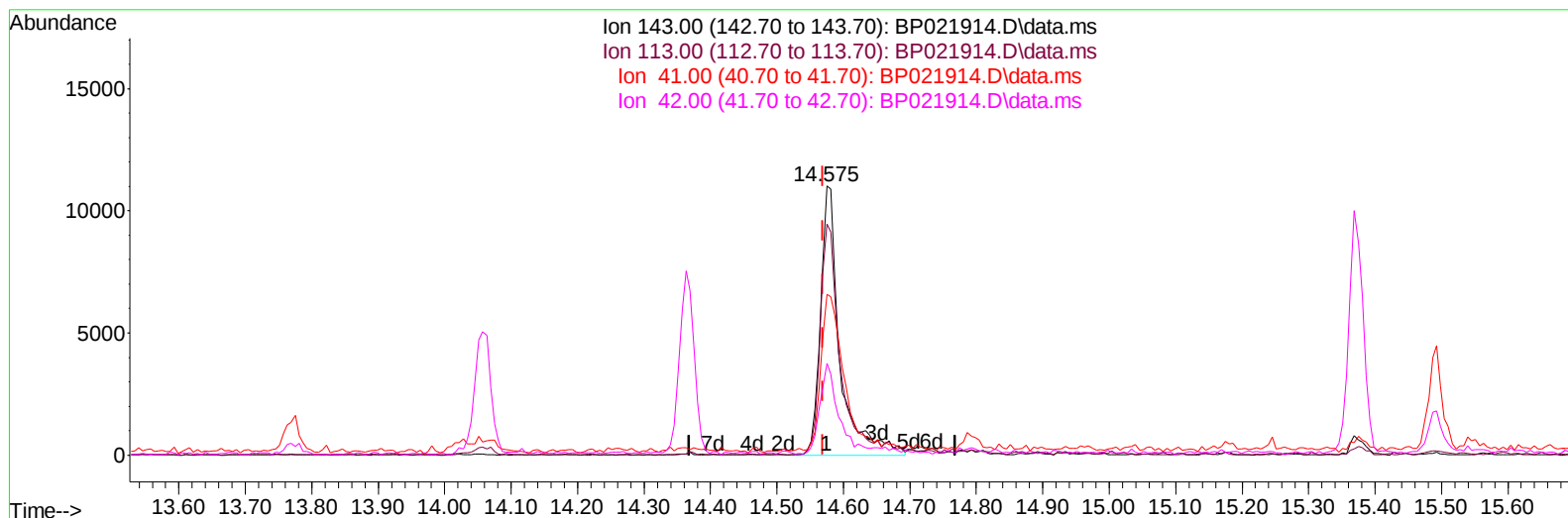
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
Data File : BP021914.D
Acq On : 13 Sep 2024 18:32
Operator : RC/JU
Sample : P3952-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B65

Manual IntegrationsAPPROVED

Quant Time: Sep 14 02:28:41 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 12 12:48:41 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/16/2024
Supervised By :mohammad ahmed 09/16/2024



TIC: BP021914.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.006) 4.66 ng/ul m

response 22356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	90.10	85.95
41.00	39.20	59.90#
42.00	26.70	34.09#

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Quant Time: Sep 14 02:30:47 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	113246	20.000	ng/ul	0.02
20) Naphthalene-d8	10.522	136	461094	20.000	ng/ul	0.02
38) Acenaphthene-d10	14.369	164	337503	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.151	188	781230	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	868211	20.000	ng/ul	-0.02
88) Perylene-d12	24.904	264	1003607	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	11836	3.381	ng/uL	0.01
4) Pyridine-d5	3.687	84	72896	7.072	ng/ul	0.02
7) Phenol-d5	6.934	99	43082	4.452	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.093	67	83856	14.920	ng/ul	0.02
11) 2-Chlorophenol-d4	7.299	132	113491	16.849	ng/ul	0.02
15) 4-Methylphenol-d8	8.452	113	125310	16.220	ng/ul	0.02
21) Nitrobenzene-d5	8.899	128	77568	27.011	ng/ul	0.02
24) 2-Nitrophenol-d4	9.616	143	80863	24.770	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.152	165	183654	25.404	ng/ul	0.02
31) 4-Chloroaniline-d4	10.652	131	240763	22.852	ng/ul	0.01
46) Dimethylphthalate-d6	13.769	166	633720	24.993	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	650011	23.370	ng/ul	0.00
54) 4-Nitrophenol-d4	14.575	143	22356m	4.664	ng/ul	0.00
60) Fluorene-d10	15.375	176	545448	22.632	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	72620	16.248	ng/ul	0.00
73) Anthracene-d10	17.251	188	960745	26.474	ng/ul	0.00
81) Pyrene-d10	19.610	212	1229810	25.602	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.668	264	1340324	25.662	ng/ul	-0.02

Target Compounds Qvalue

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

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