

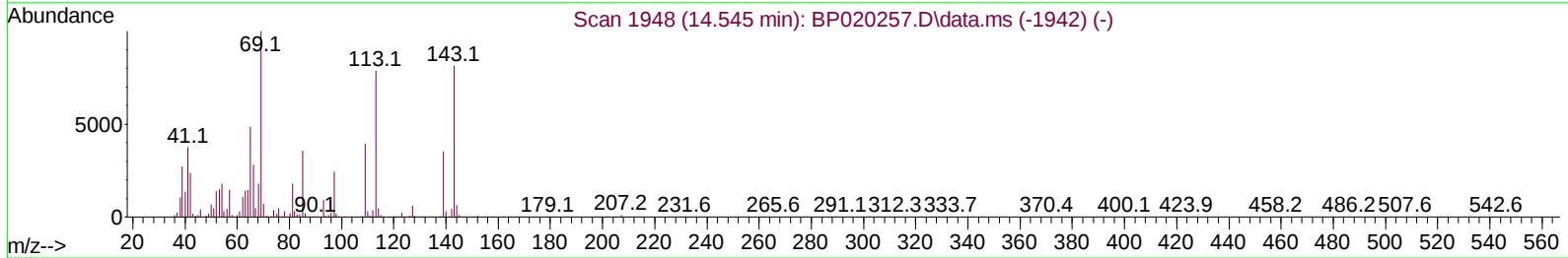
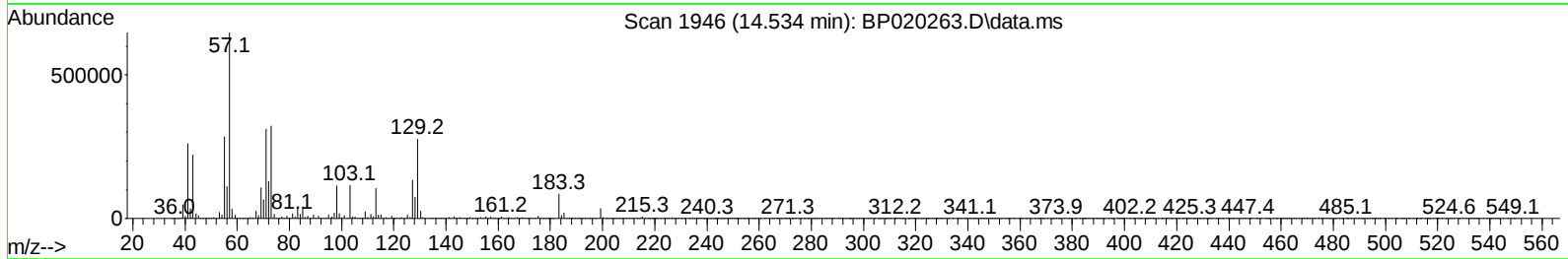
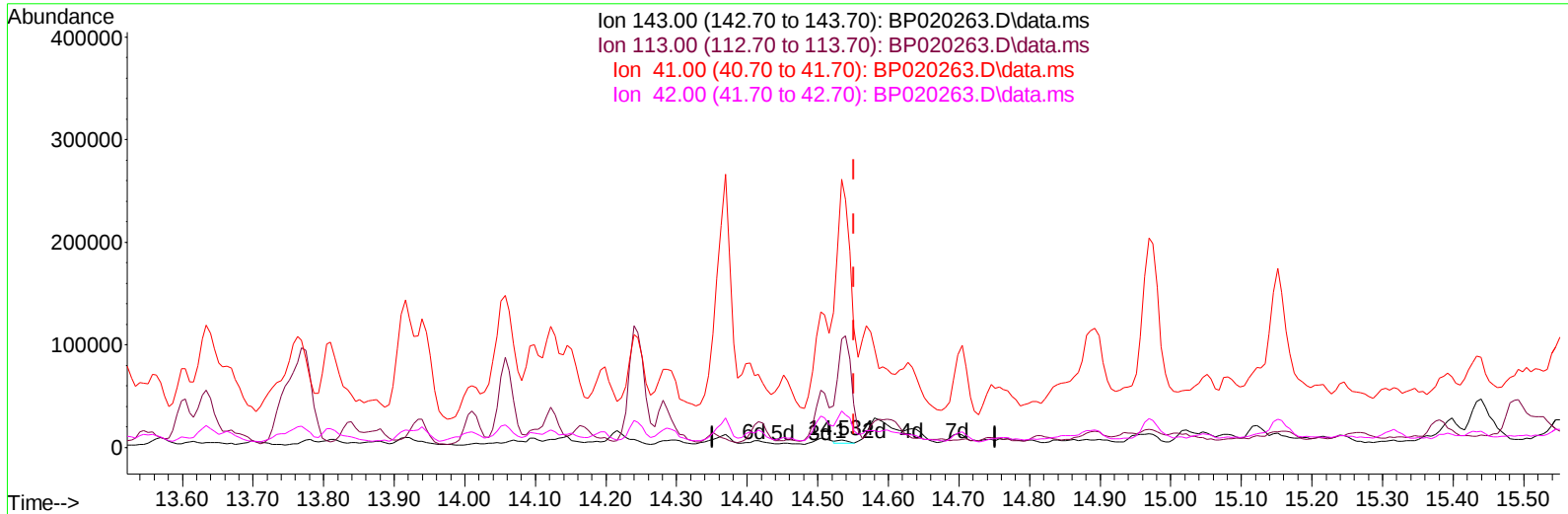
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020263.D
Acq On : 09 May 2024 10:00
Operator : MA/JU
Sample : P2369-13
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N7

Manual IntegrationsAPPROVED

Quant Time: May 09 10:57:56 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 08 22:57:55 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/09/2024
Supervised By :mohammad ahmed 05/11/2024



TIC: BP020263.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.534min (-0.018) 1.27 ng/u1

response 3186

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	1455.04#
41.00	50.30	3597.30#
42.00	32.50	484.83#

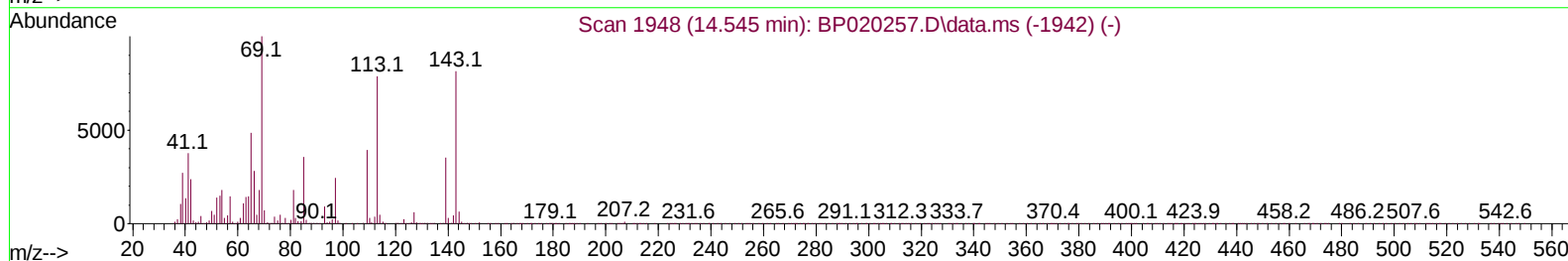
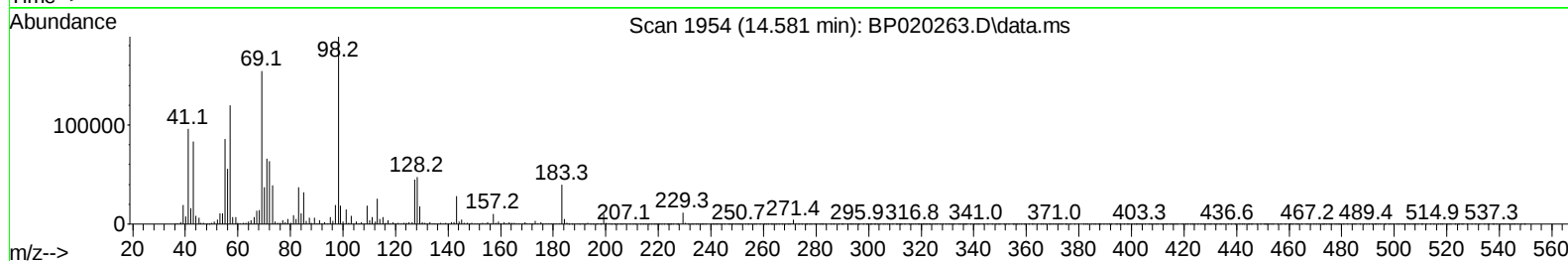
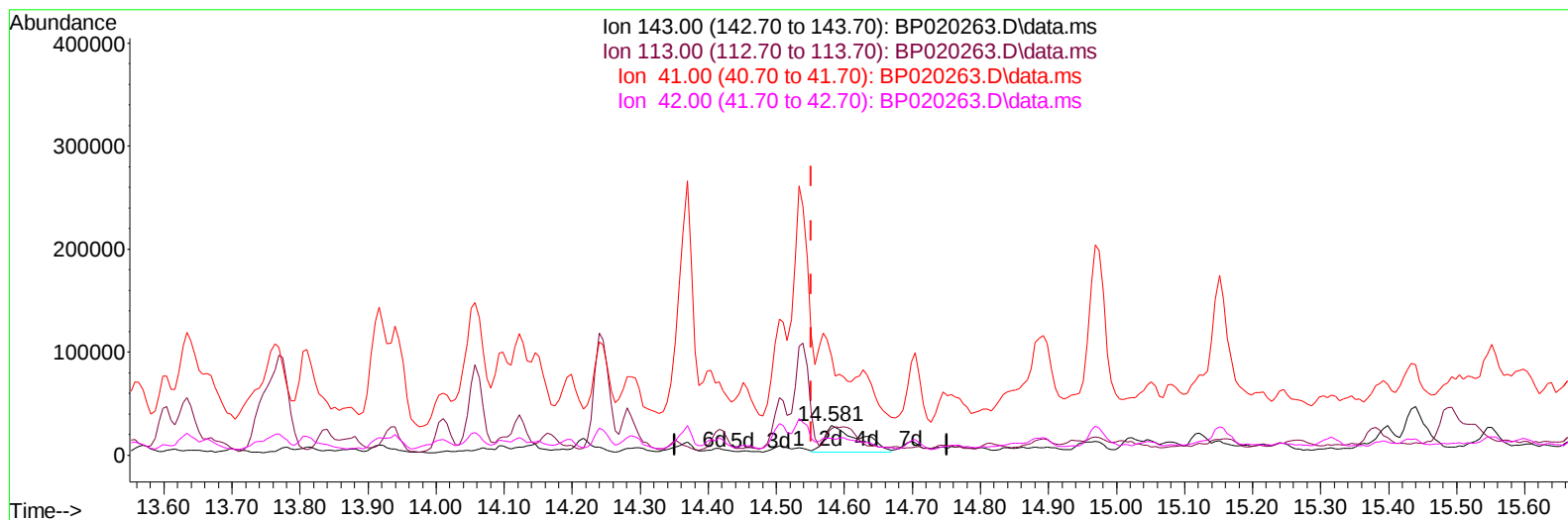
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(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.029) 36.74 ng/ul m

response 91968

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	89.94
41.00	50.30	338.41#
42.00	32.50	57.39#

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/09/2024
 Supervised By :mohammad ahmed 05/11/2024

Quant Time: May 09 11:05:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.640	152		99041	20.000	ng/ul	0.02
20) Naphthalene-d8	10.416	136		374582	20.000	ng/ul	# 0.03
38) Acenaphthene-d10	14.298	164		223009	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.151	188		442027	20.000	ng/ul	0.01
79) Chrysene-d12	21.639	240		484428	20.000	ng/ul	0.00
88) Perylene-d12	25.009	264		580925	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.222	96		15208	6.485	ng/uL	0.01
4) Pyridine-d5	3.628	84		180455	26.483	ng/ul	0.01
7) Phenol-d5	6.822	99		355780	41.829	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.010	67		787275	152.371	ng/ul	0.04
11) 2-Chlorophenol-d4	7.175	132		184729	30.381	ng/ul	0.02
15) 4-Methylphenol-d8	8.322	113		648755	94.343	ng/ul	0.00
21) Nitrobenzene-d5	8.804	128		104969	37.422	ng/ul	0.02
24) 2-Nitrophenol-d4	9.516	143		98906	30.791	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.057	165		175020	29.858	ng/ul	0.04
31) 4-Chloroaniline-d4	10.569	131		93381	11.383	ng/ul	0.02
46) Dimethylphthalate-d6	13.710	166		451040	27.696	ng/ul	0.00
49) Acenaphthylene-d8	13.987	160		520105	28.911	ng/ul	0.00
54) 4-Nitrophenol-d4	14.581	143		91968m	36.744	ng/ul	0.03
60) Fluorene-d10	15.316	176		379005	27.902	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200		50031	17.836	ng/ul	0.00
73) Anthracene-d10	17.251	188		584077	30.544	ng/ul	0.00
81) Pyrene-d10	19.645	212		739891	26.129	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.786	264		812803	28.937	ng/ul	0.01
Target Compounds							
71) Pentachlorophenol	16.845	266		16539639	4934.394	ng/ul	Qvalue 98

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

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