

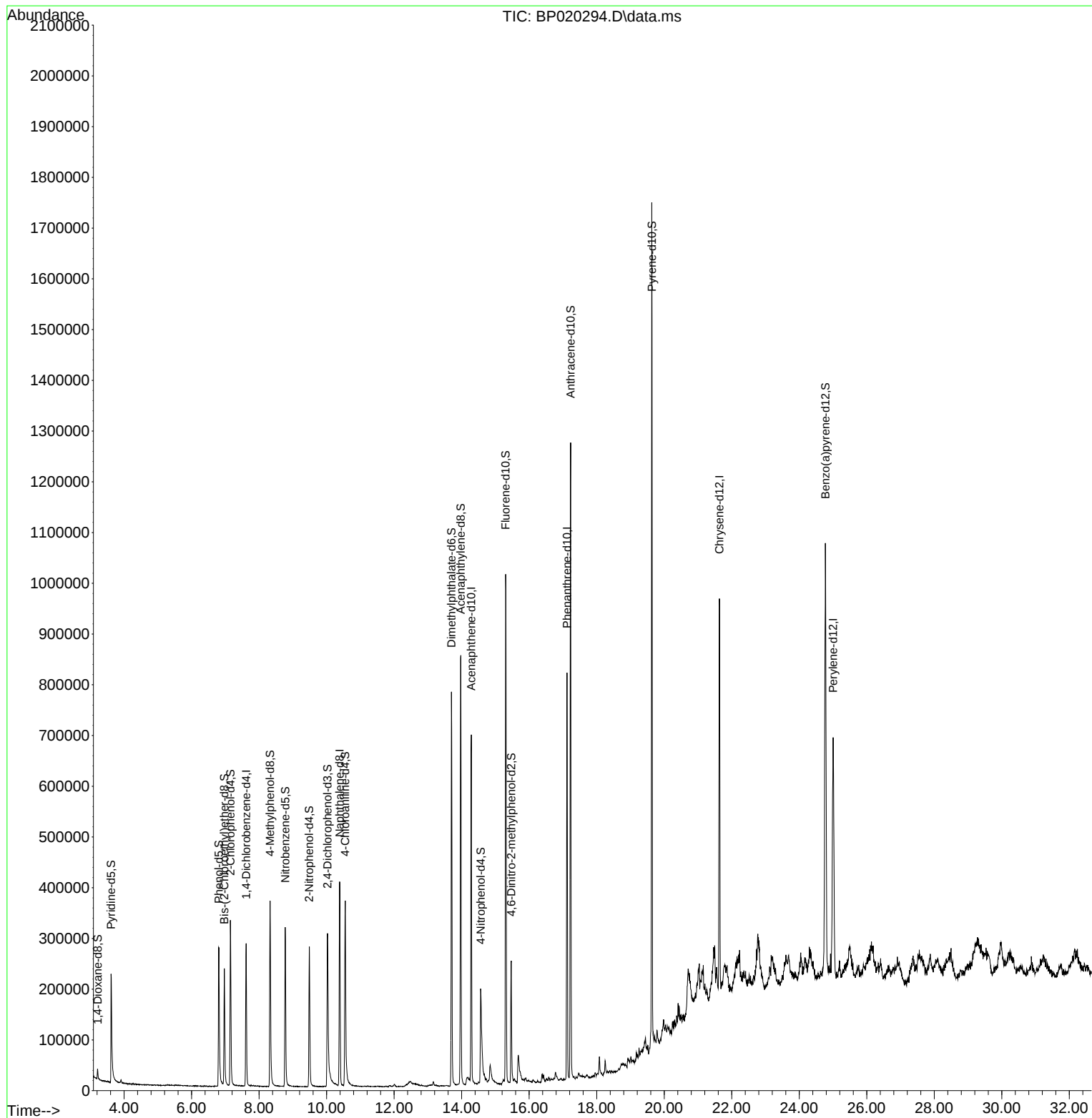
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020294.D
Acq On : 10 May 2024 17:09
Operator : MA/JU
Sample : PB160761BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK761

Manual IntegrationsAPPROVED

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

Quant Time: May 10 23:56:08 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



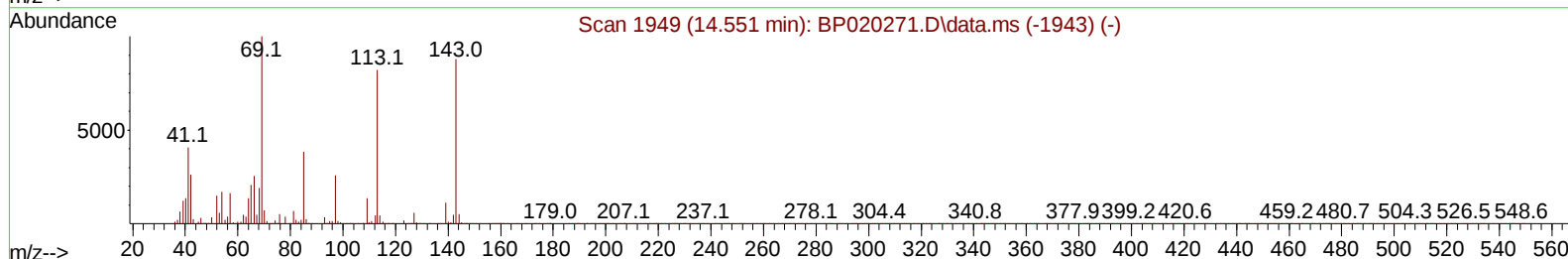
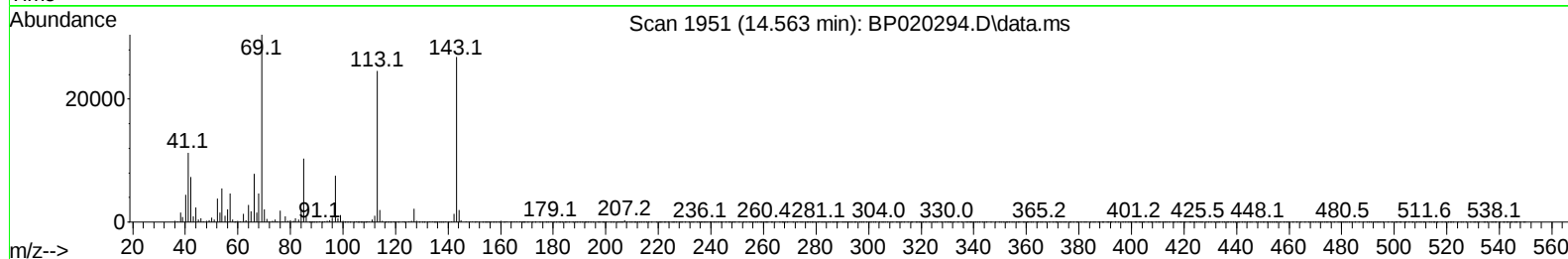
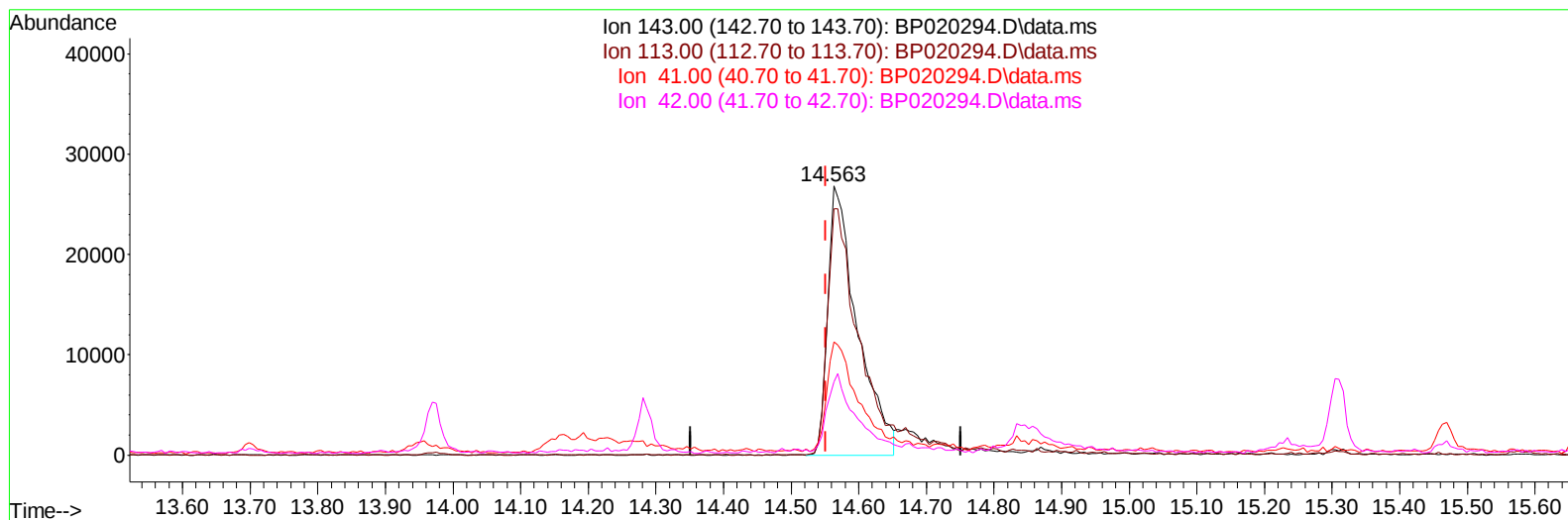
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TIC: BP020294.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 29.67 ng/u1

response 81081

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	91.56
41.00	50.30	42.02
42.00	32.50	27.20

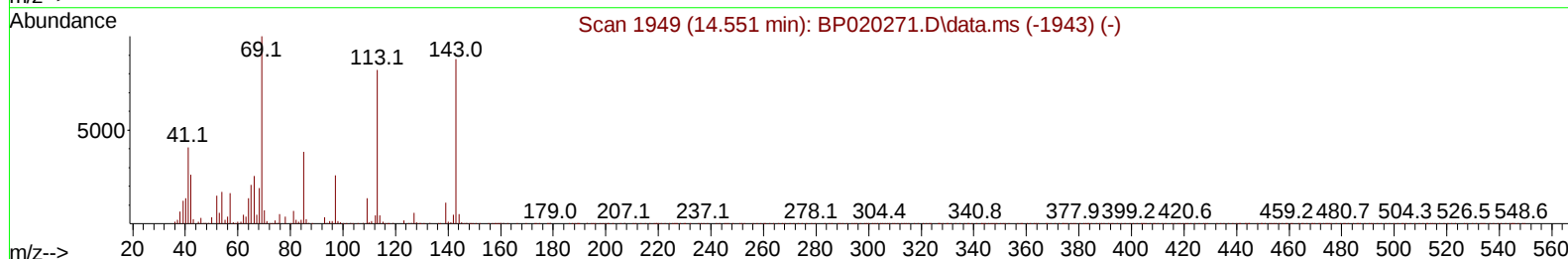
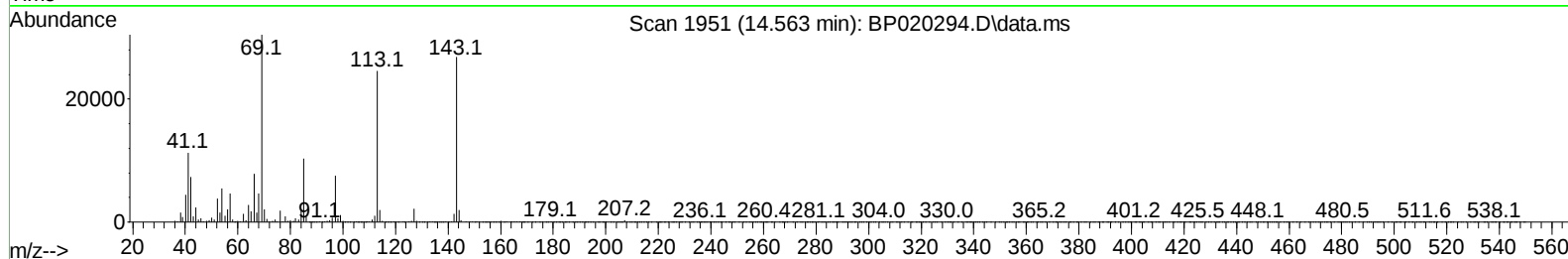
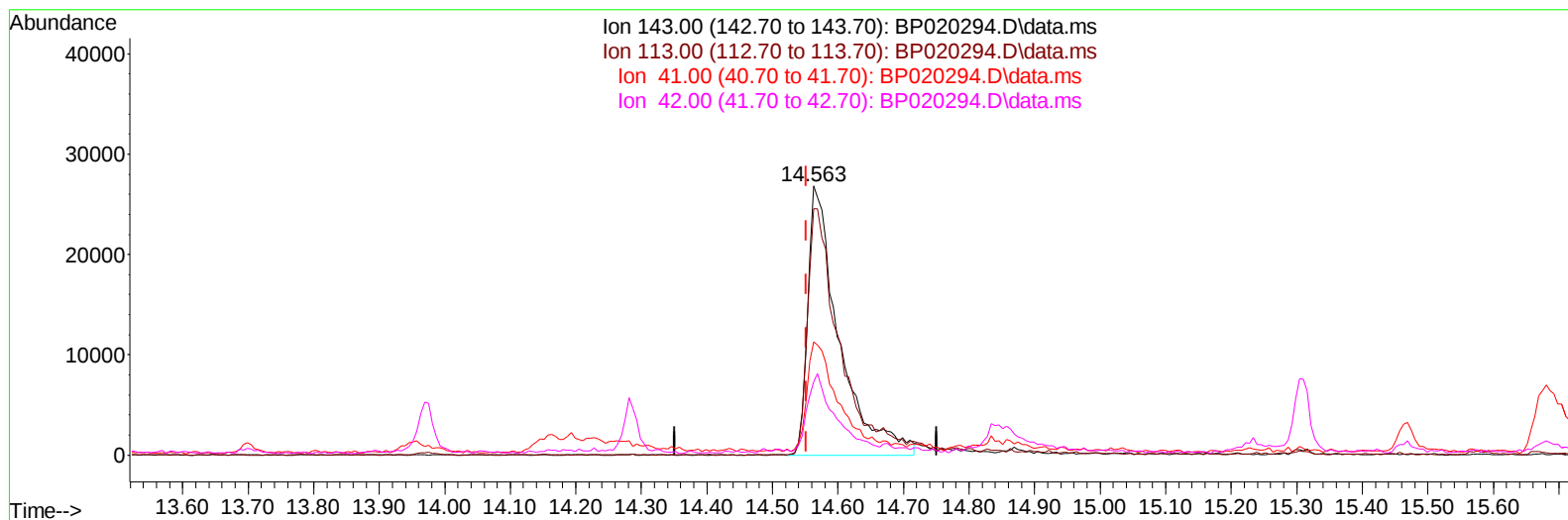
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TIC: BP020294.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 32.40 ng/ul m

response 88536

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	91.56
41.00	50.30	42.02
42.00	32.50	27.20

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 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.617	152	82918	20.000	ng/u1	0.00
20) Naphthalene-d8	10.387	136	343417	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.287	164	243496	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.122	188	521049	20.000	ng/u1	-0.02
79) Chrysene-d12	21.633	240	499896	20.000	ng/u1	0.00
88) Perylene-d12	25.004	264	574491	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	10496	5.346	ng/uL	0.00
4) Pyridine-d5	3.617	84	143382	25.134	ng/u1	0.00
7) Phenol-d5	6.811	99	191964	26.958	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	113158	26.159	ng/u1	0.00
11) 2-Chlorophenol-d4	7.152	132	151106	29.684	ng/u1	0.00
15) 4-Methylphenol-d8	8.328	113	158082	27.459	ng/u1	0.00
21) Nitrobenzene-d5	8.775	128	77632	30.188	ng/u1	0.00
24) 2-Nitrophenol-d4	9.487	143	89496	30.390	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	157202	29.253	ng/u1	0.00
31) 4-Chloroaniline-d4	10.552	131	230403	30.633	ng/u1	0.00
46) Dimethylphthalate-d6	13.699	166	535750	30.130	ng/u1	-0.01
49) Acenaphthylene-d8	13.969	160	578580	29.455	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.563	143	88536m	32.396	ng/u1	0.01
60) Fluorene-d10	15.304	176	451280	30.428	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.469	200	61827	18.699	ng/u1	-0.01
73) Anthracene-d10	17.228	188	752385	33.379	ng/u1	-0.02
81) Pyrene-d10	19.634	212	904486	30.953	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.774	264	906974	32.652	ng/u1	0.00

Target Compounds Qvalue

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