

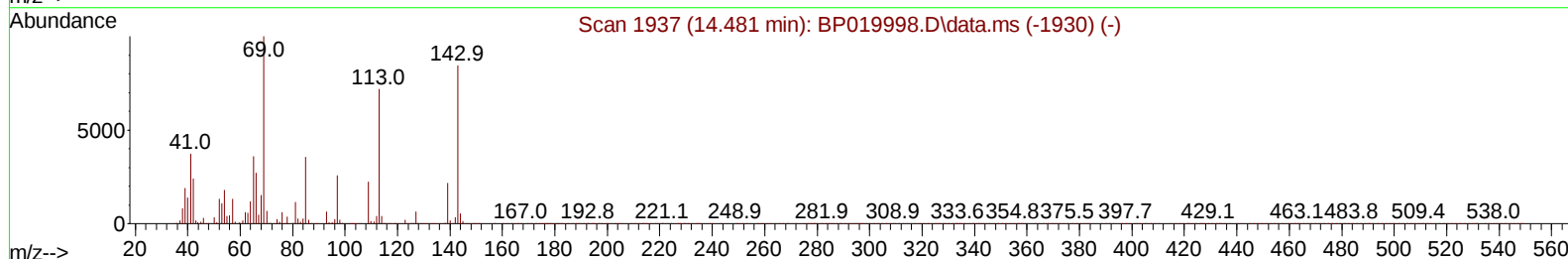
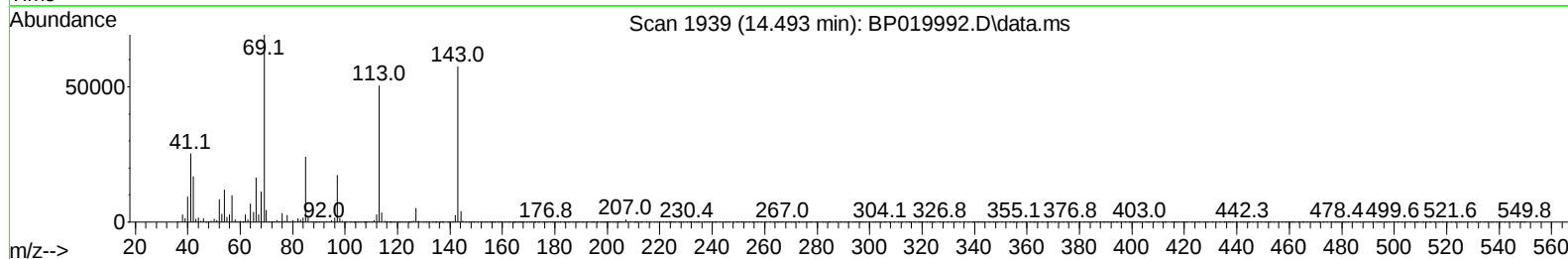
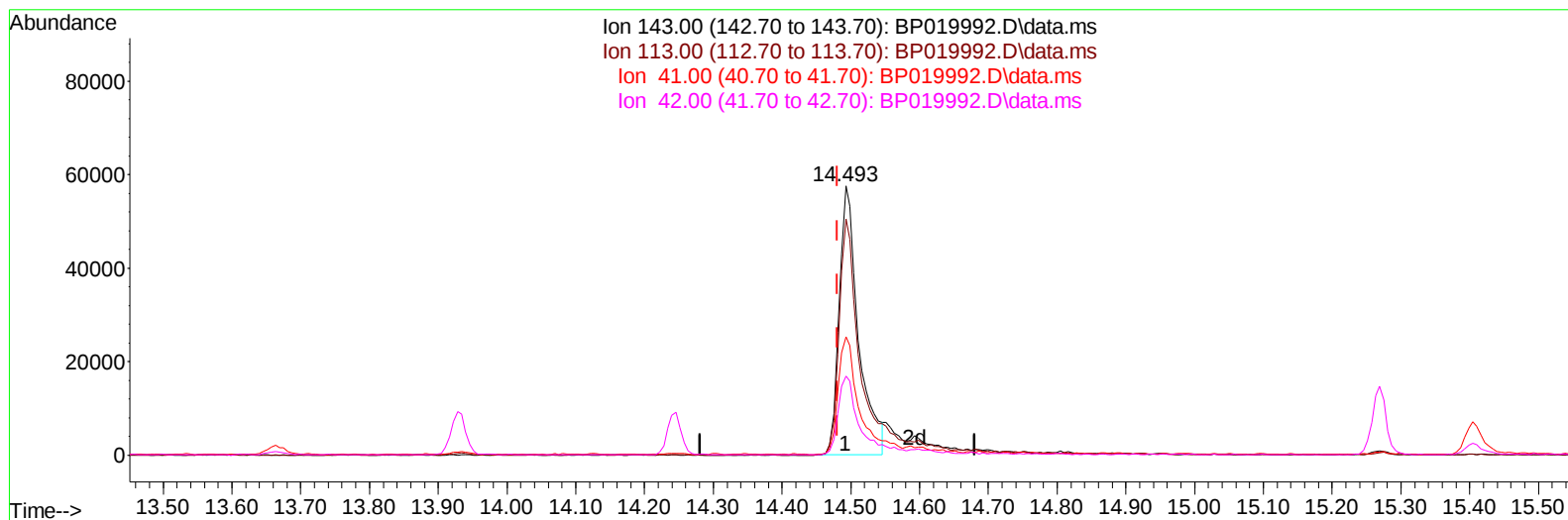
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032624\
 Data File : BP019992.D
 Acq On : 26 Mar 2024 10:24
 Operator : MA/JU
 Sample : PB159439BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK439

Manual IntegrationsAPPROVED

Quant Time: Mar 26 11:19:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :mohammad ahmed 03/29/2024



TIC: BP019992.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.012) 26.39 ng/u1

response 112994

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	87.83
41.00	40.90	43.99
42.00	28.50	29.37

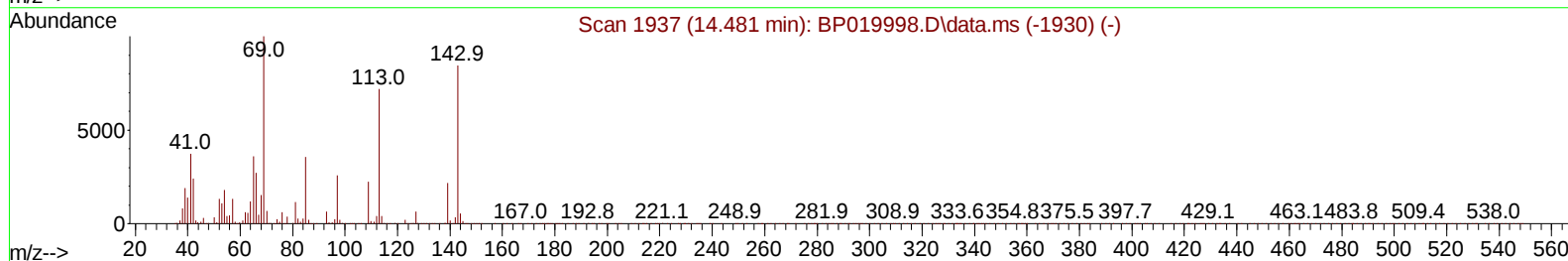
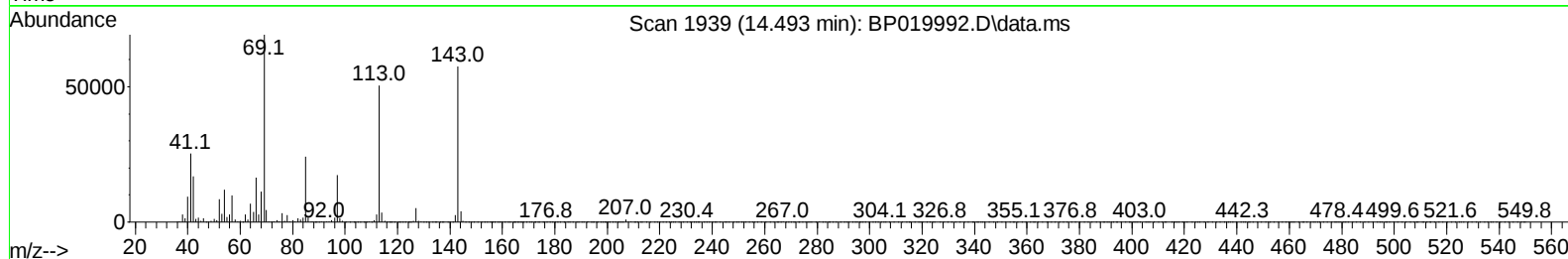
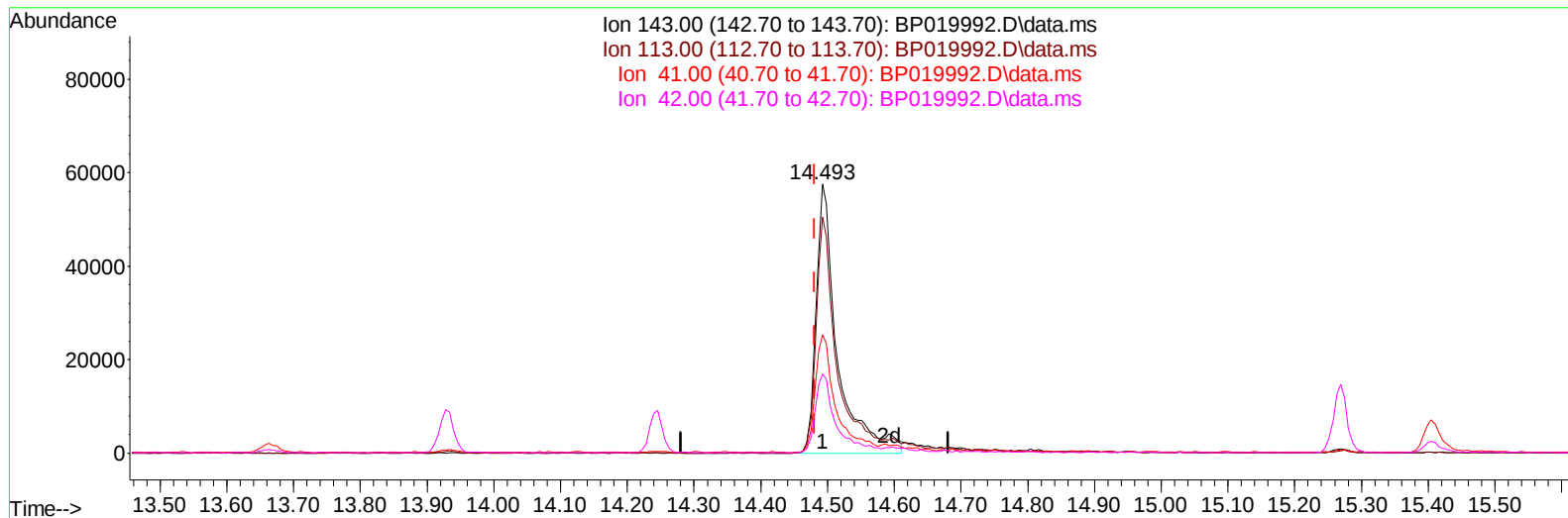
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TIC: BP019992.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.012) 30.05 ng/ul m

response 128661

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	87.83
41.00	40.90	43.99
42.00	28.50	29.37

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 Misc :
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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :mohammad ahmed 03/29/2024

Quant Time: Mar 26 23:51:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	138850	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	583609	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.246	164	388418	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.063	188	790026	20.000	ng/ul	-0.01
79) Chrysene-d12	21.533	240	612388	20.000	ng/ul	-0.01
88) Perylene-d12	24.816	264	639065	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	21494	6.783	ng/uL	0.00
4) Pyridine-d5	3.640	84	308131	32.201	ng/ul	0.00
7) Phenol-d5	6.811	99	388899	33.210	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.969	67	275342	36.089	ng/ul	0.00
11) 2-Chlorophenol-d4	7.164	132	284731	33.856	ng/ul	0.01
15) 4-Methylphenol-d8	8.322	113	316211	34.901	ng/ul	0.02
21) Nitrobenzene-d5	8.764	128	152952	34.193	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	172871	34.540	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	285533	31.238	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	413498	34.702	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	976164	35.845	ng/ul	0.01
49) Acenaphthylene-d8	13.934	160	1096923	34.481	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	128661m	30.052	ng/ul	0.01
60) Fluorene-d10	15.269	176	818817	35.296	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	140545	29.720	ng/ul	-0.02
73) Anthracene-d10	17.181	188	1197746	34.249	ng/ul	0.01
81) Pyrene-d10	19.563	212	1318651	36.701	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	1095507	35.682	ng/ul	-0.02

Target Compounds Qvalue

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

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