

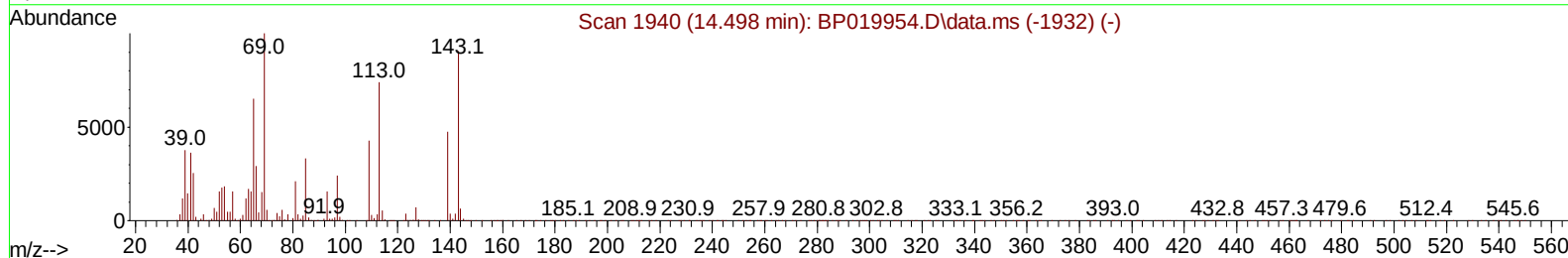
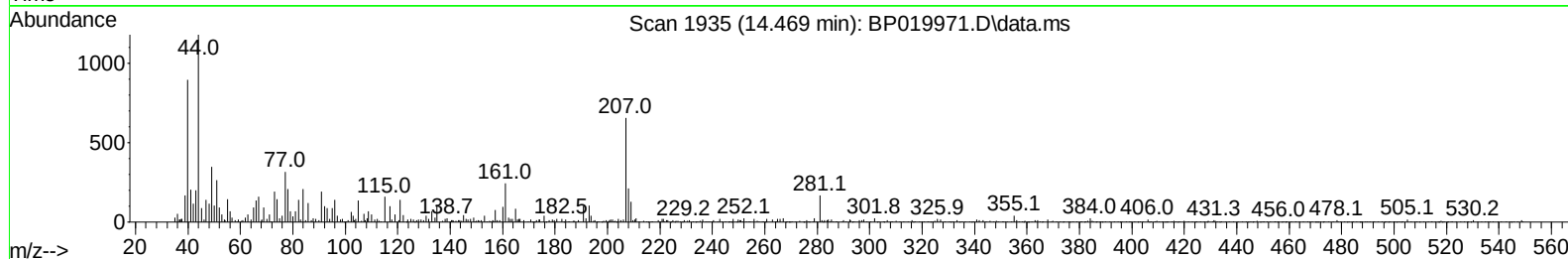
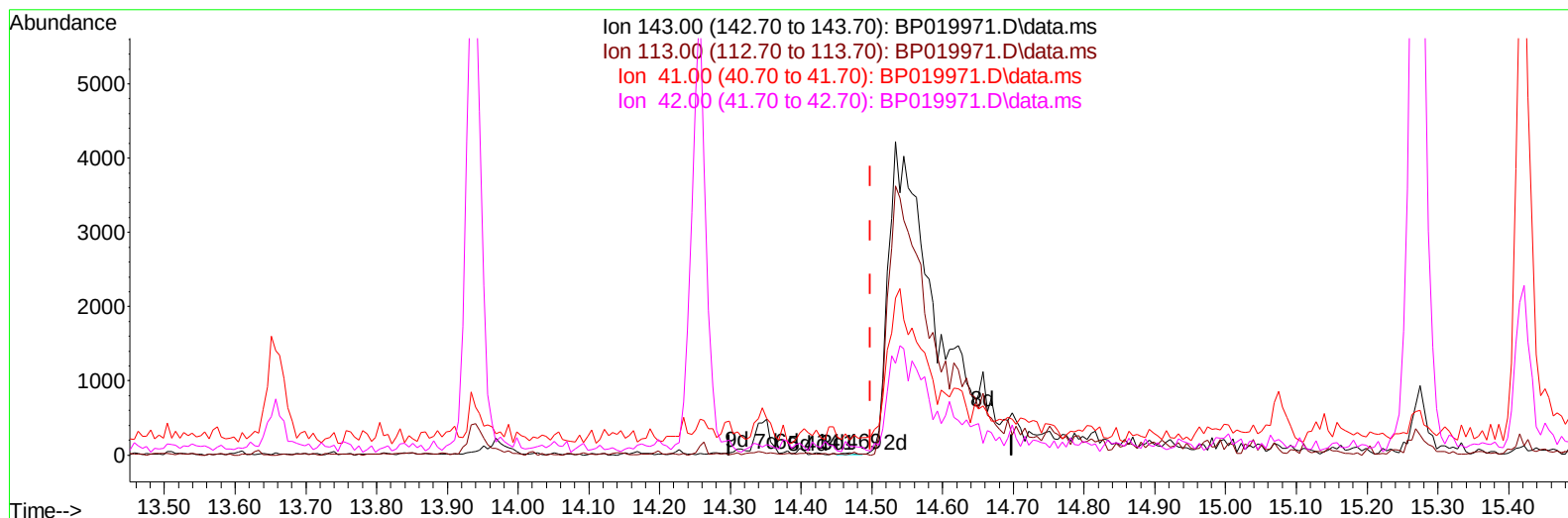
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
 Data File : BP019971.D
 Acq On : 23 Mar 2024 01:39
 Operator : MA/JU
 Sample : P1841-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AK7

Manual IntegrationsAPPROVED

Quant Time: Mar 23 02:08:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 22 22:40:53 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/26/2024
 Supervised By :mohammad ahmed 03/26/2024



TIC: BP019971.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (-0.029) 0.01 ng/u1

response 27

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	55.56#
41.00	40.90	748.15#
42.00	28.50	429.63#

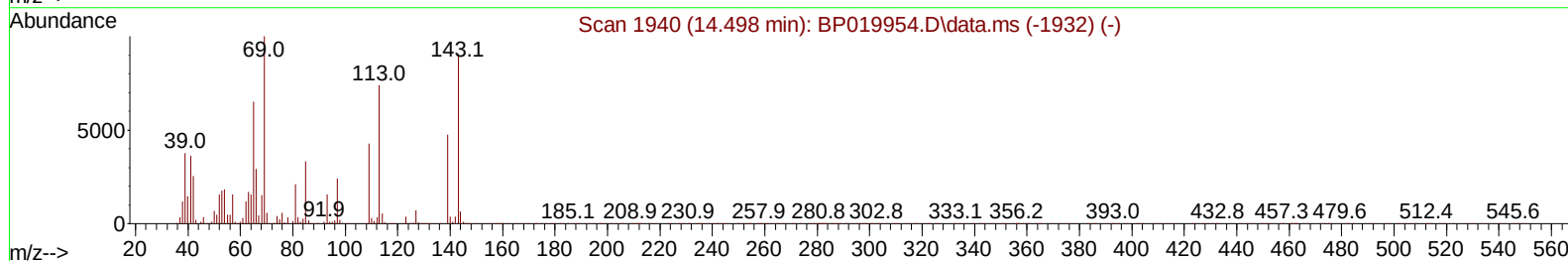
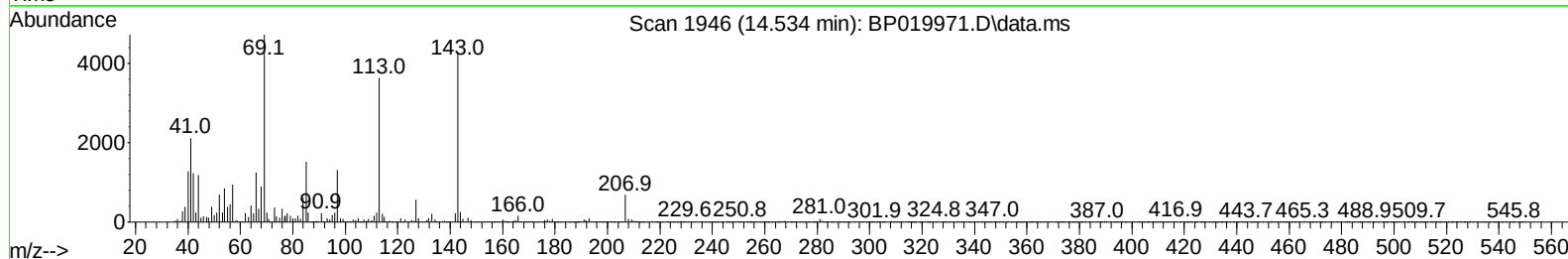
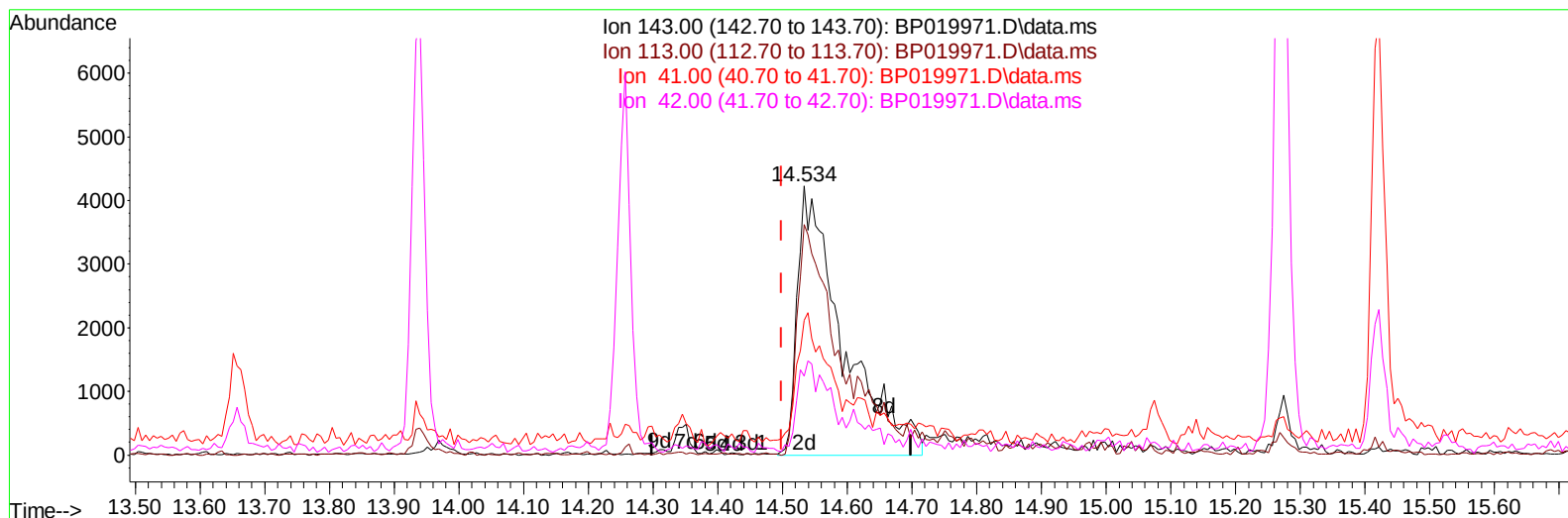
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(54) 4-Nitrophenol-d4 (S)

14.534min (+ 0.035) 6.82 ng/ul m

response 20568

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	85.71
41.00	40.90	50.07#
42.00	28.50	29.29

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.622	152		101025	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136		413023	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164		273547	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188		577303	20.000	ng/ul	-0.02
79) Chrysene-d12	21.551	240		491817	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264		540920	20.000	ng/ul	-0.04
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.246	96		11994	5.202	ng/uL	0.00
4) Pyridine-d5	3.652	84		43554	6.256	ng/ul	0.00
7) Phenol-d5	6.822	99		56903	6.678	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.975	67		198892	35.829	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132		162557	26.566	ng/ul	0.01
15) 4-Methylphenol-d8	8.322	113		110520	16.766	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128		112809	35.634	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143		124500	35.150	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165		195562	30.232	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131		243403	28.864	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166		733328	38.236	ng/ul	-0.01
49) Acenaphthylene-d8	13.939	160		817988	36.511	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.534	143		20568m	6.822	ng/ul	0.04
60) Fluorene-d10	15.275	176		622528	38.104	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200		115875	33.532	ng/ul	-0.01
73) Anthracene-d10	17.175	188		1018592	39.859	ng/ul	-0.01
81) Pyrene-d10	19.569	212		1165851	40.403	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.604	264		1070239	41.184	ng/ul	-0.03
Target Compounds							
						Qvalue	
2) 1,4-Dioxane	3.275	88		3050	1.159	ng/ul#	67
16) Acetophenone	8.440	105		147266	13.850	ng/ul	99

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

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