

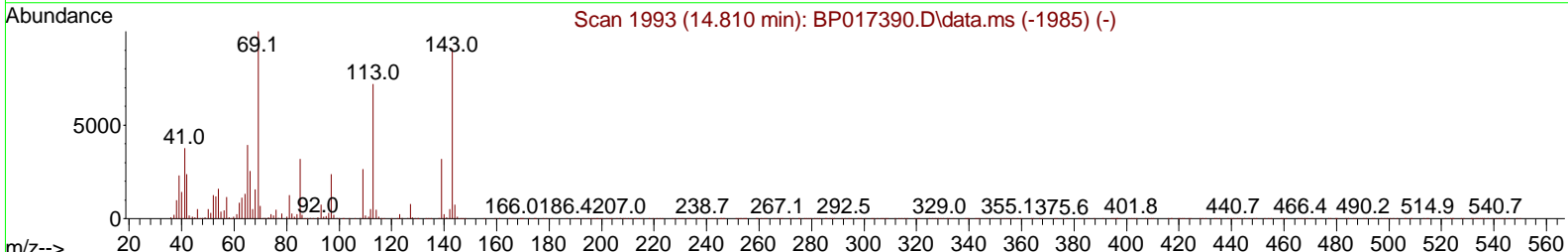
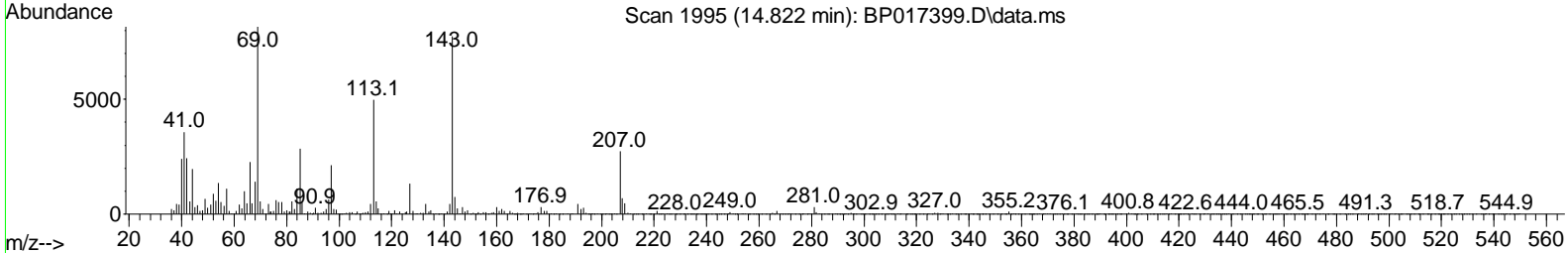
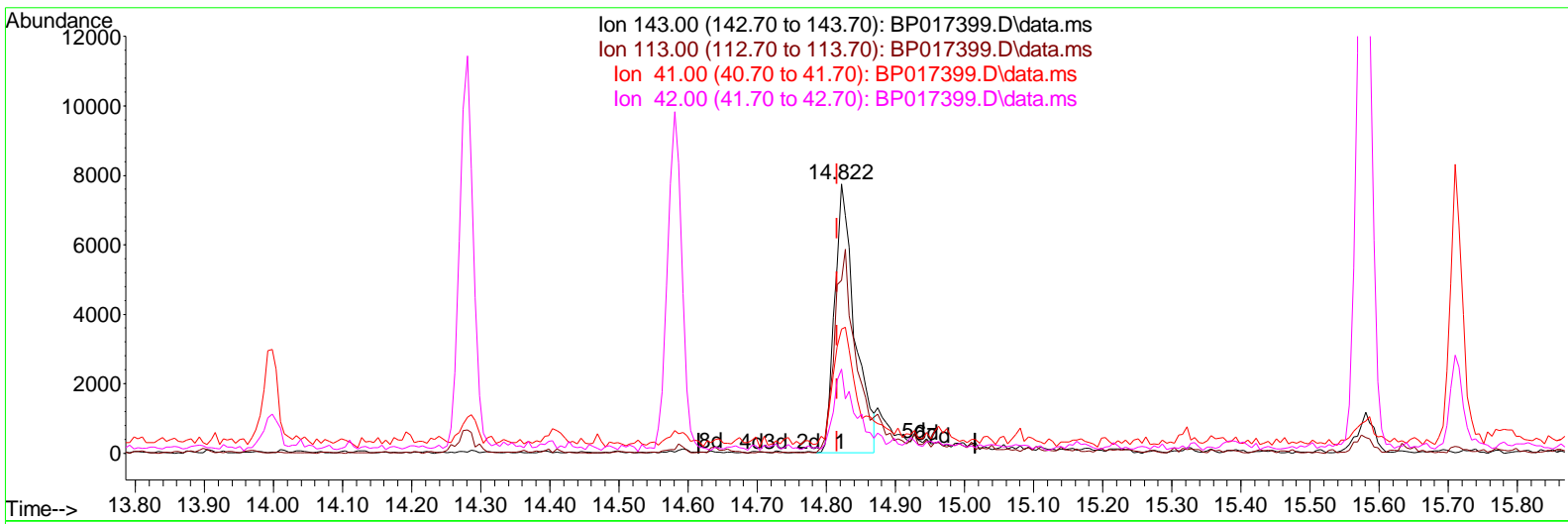
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
 Data File : BP017399.D
 Acq On : 27 Sep 2023 16:01
 Operator : MA/JU
 Sample : 04359-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 JYCA2

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/28/2023
 Supervised By :mohammad ahmed 09/29/2023

Quant Time: Sep 27 22:30:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration



TIC: BP017399.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.006) 3.09 ng/ul

response	15708	
Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	64.14
41.00	36.70	46.20#
42.00	26.40	31.32

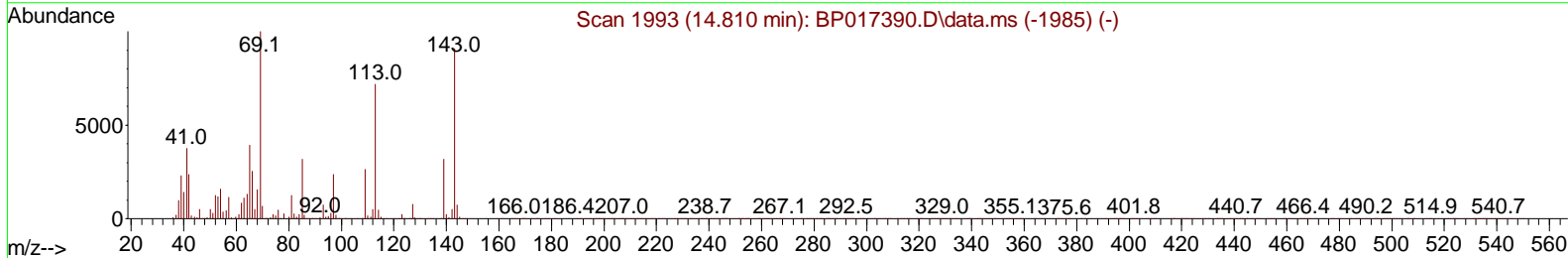
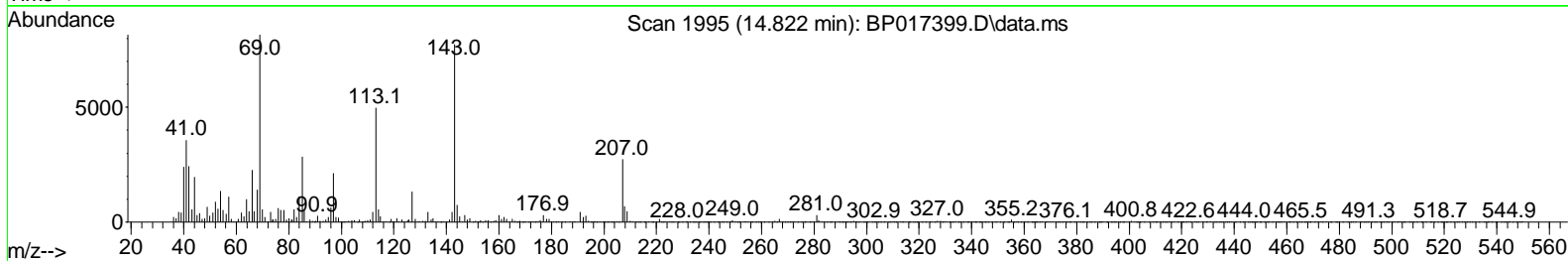
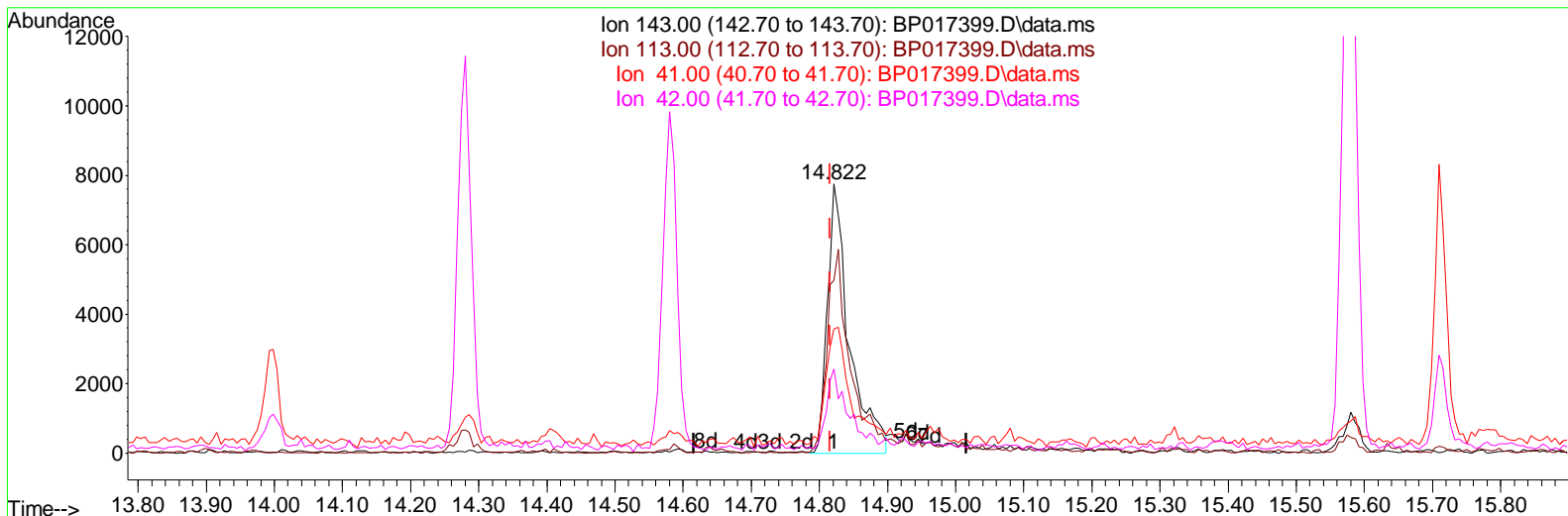
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TIC: BP017399.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.006) 3.40 ng/ul m

response 17290

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	64.14
41.00	36.70	46.20#
42.00	26.40	31.32

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
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Instrument :
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Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/28/2023
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Quant Time: Sep 27 22:57:51 2023
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Compound	R.T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	149022	20.000	ng/ul	0.00
20) Naphthalene-d8	10.739	136	661954	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.580	164	407736	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	912232	20.000	ng/ul	0.00
79) Chrysene-d12	21.456	240	839280	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	938872	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	8773	2.418	ng/uL	0.00
4) Pyridine-d5	3.728	84	115674	11.405	ng/ul	0.00
7) Phenol-d5	7.075	99	75277	6.151	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	271688	36.783	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	233215	24.025	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	148465	15.603	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	164601	34.396	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	155683	29.961	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	268588	27.359	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	410610	28.874	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	1150962	39.404	ng/ul	0.00
49) Acenaphthylene-d8	14.280	160	1217941	36.255	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	17290m	3.398	ng/ul	0.00
60) Fluorene-d10	15.580	176	1014157	40.330	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	119415	23.226	ng/ul	0.00
73) Anthracene-d10	17.457	188	1743880	43.103	ng/ul	0.00
81) Pyrene-d10	19.698	212	2068973	45.586	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.809	264	1999818	44.240	ng/ul	0.00

Target Compounds Qvalue

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

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Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION
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