

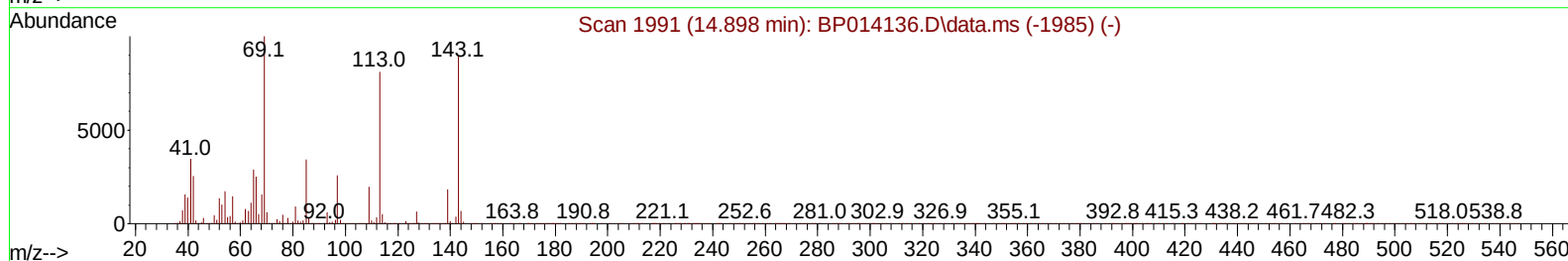
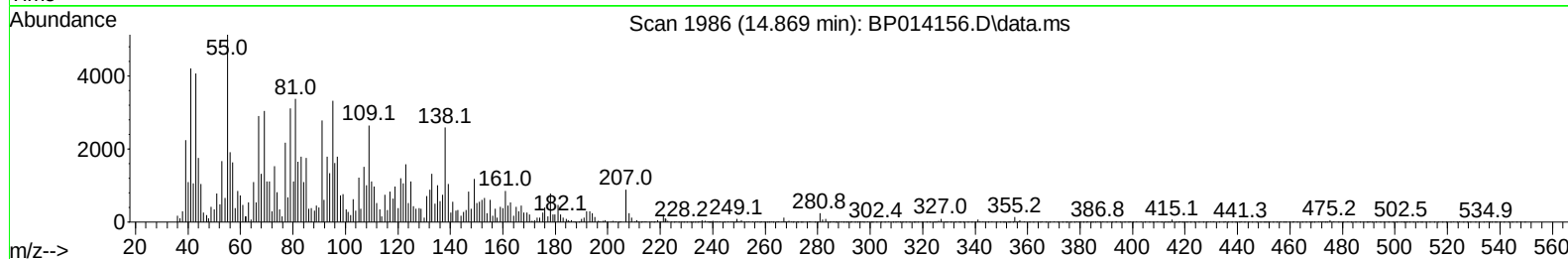
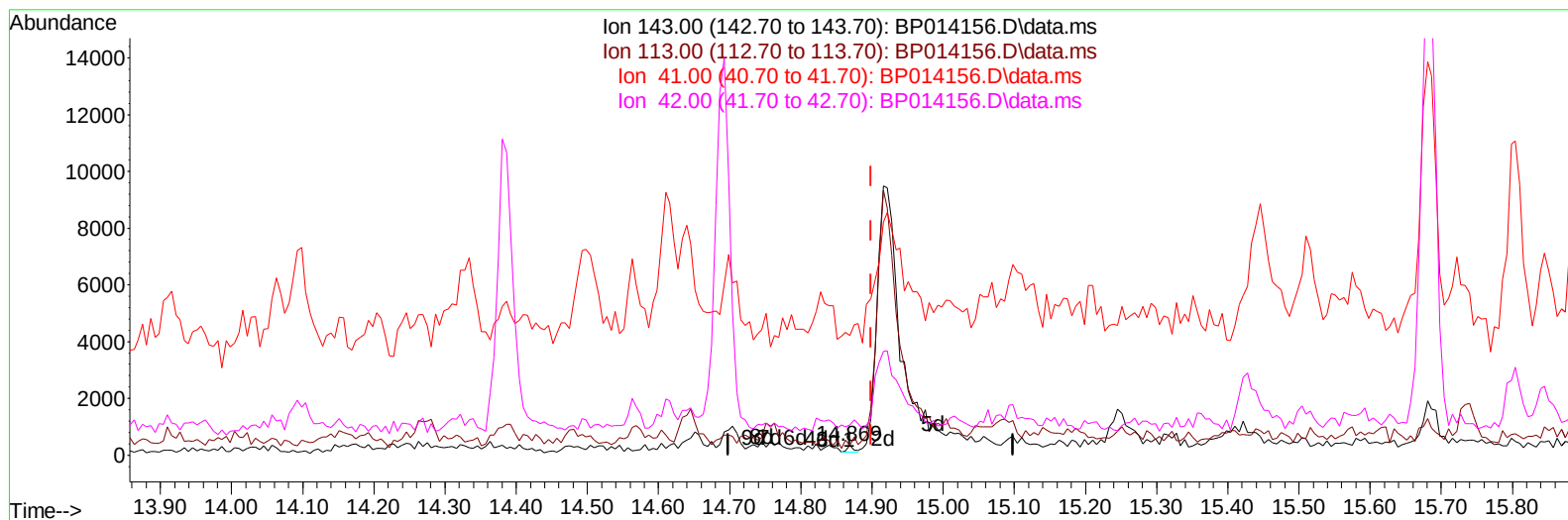
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032023\
Data File : BP014156.D
Acq On : 21 Mar 2023 00:10
Operator : CG/JU
Sample : 01885-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB908

Quant Time: Mar 21 00:50:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 20 13:12:59 2023
Response via : Initial Calibration

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/21/2023
Supervised By :Jagrut Upadhyay 03/21/2023



TIC: BP014156.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.869min (-0.029) 0.02 ng/u1

response 131

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	101.77
41.00	40.20	1243.07#
42.00	28.50	312.68#

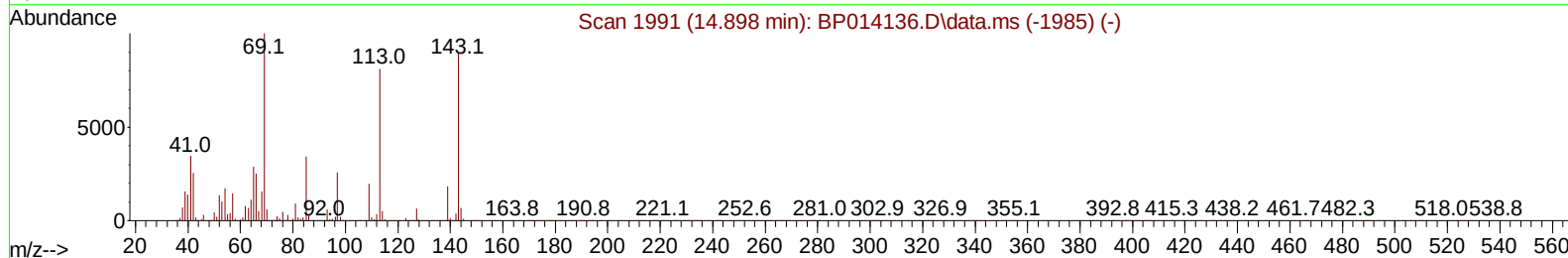
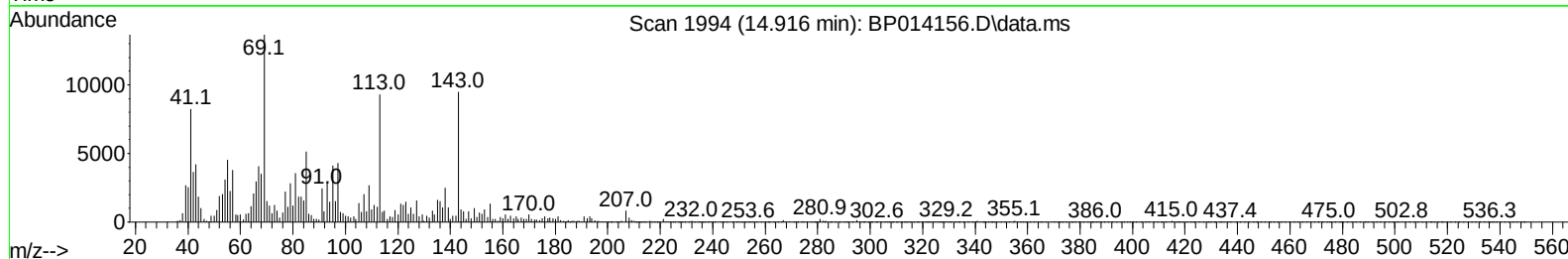
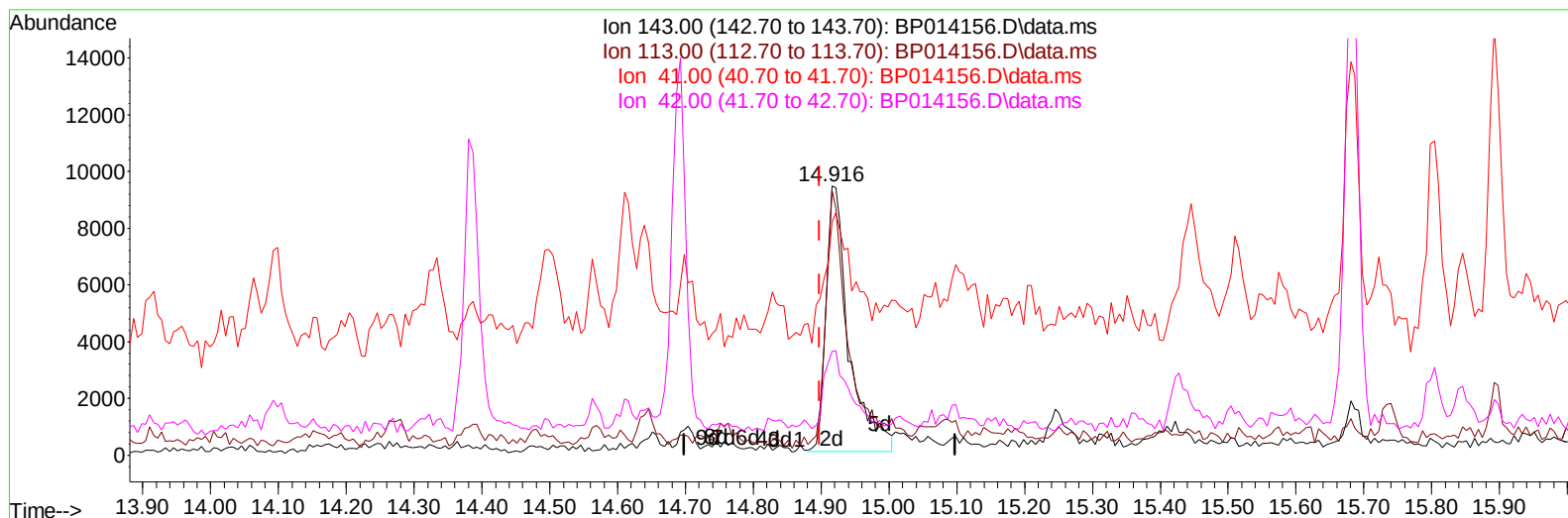
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(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.018) 2.55 ng/ul m

response 22095

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	98.35
41.00	40.20	86.67#
42.00	28.50	38.52#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.052	152	200093	20.000	ng/ul	0.00
20) Naphthalene-d8	10.875	136	844359	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.692	164	579511	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	1315823	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.521	240	1228956	20.000	ng/ul	#-0.01
88) Perylene-d12	24.045	264	1119828	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	15325	2.894	ng/uL	0.00
4) Pyridine-d5	3.852	84	140594	9.588	ng/ul	0.00
7) Phenol-d5	7.210	99	64370	3.645	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	241327	23.434	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132	150433	11.182	ng/ul	0.00
15) 4-Methylphenol-d8	8.763	113	141736	9.827	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	153494	22.783	ng/ul	0.00
24) 2-Nitrophenol-d4	9.951	143	111493	13.825	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.493	165	203905	13.437	ng/ul	0.00
31) 4-Chloroaniline-d4	11.016	131	314081	15.640	ng/ul	0.00
46) Dimethylphthalate-d6	14.092	166	1170573	23.801	ng/ul	0.00
49) Acenaphthylene-d8	14.386	160	1178104	22.612	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	22095m	2.549	ng/ul	0.02
60) Fluorene-d10	15.681	176	948964	22.767	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.804	200	108909	10.951	ng/ul	0.00
73) Anthracene-d10	17.545	188	1561963	24.453	ng/ul	-0.01
81) Pyrene-d10	19.769	212	1931743	28.500	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	1490852	25.028	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.452	88	127245	22.257	ng/uL	92

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

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