

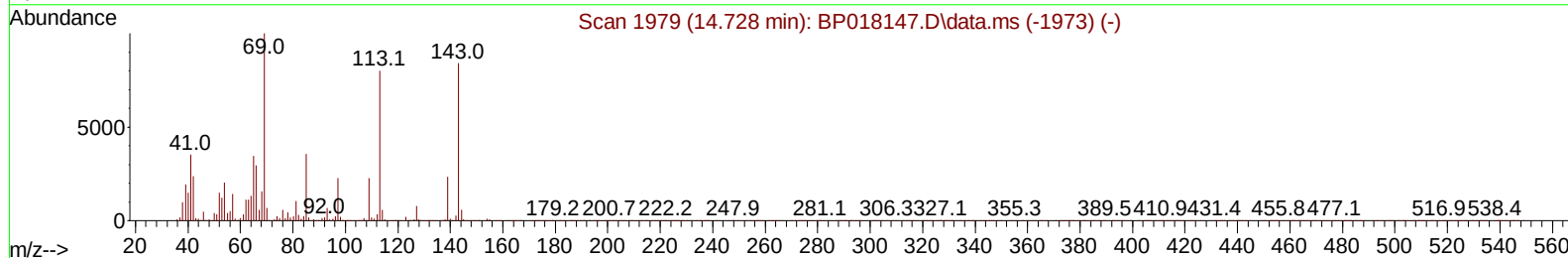
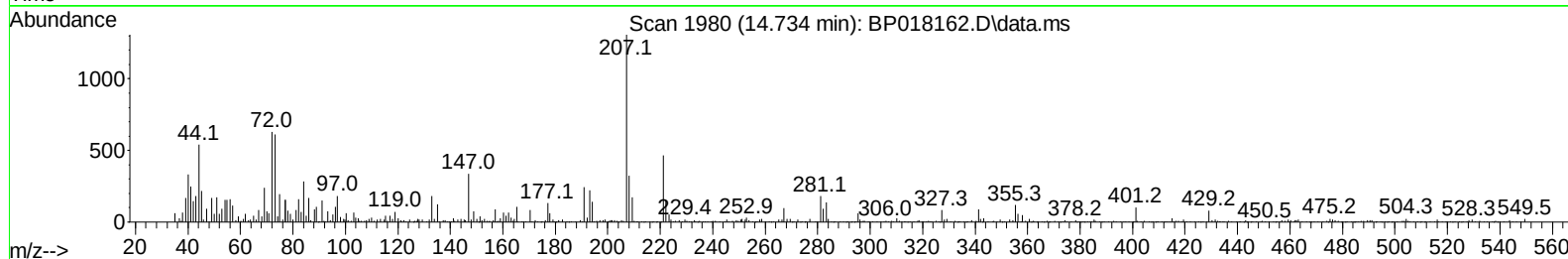
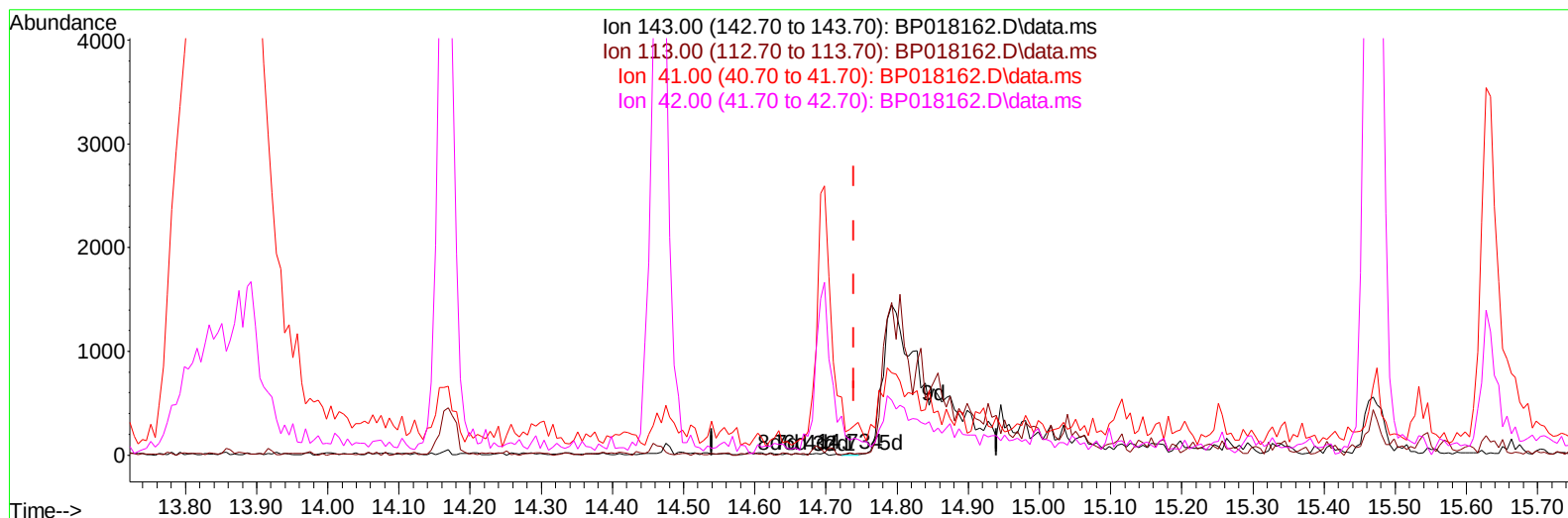
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018162.D
Acq On : 14 Nov 2023 18:58
Operator : MA/JU
Sample : 05337-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDJ3

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:39:58 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023



TIC: BP018162.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 0.00 ng/u1

response 12

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	150.00#
41.00	47.70	1556.25#
42.00	30.10	918.75#

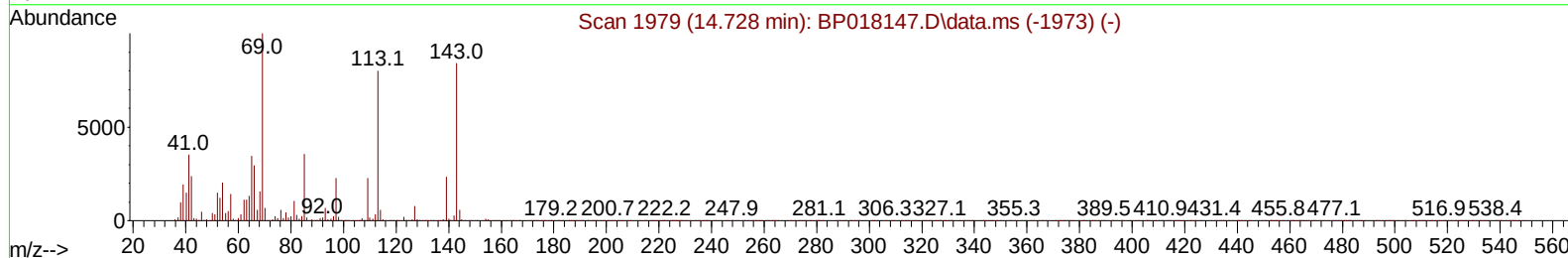
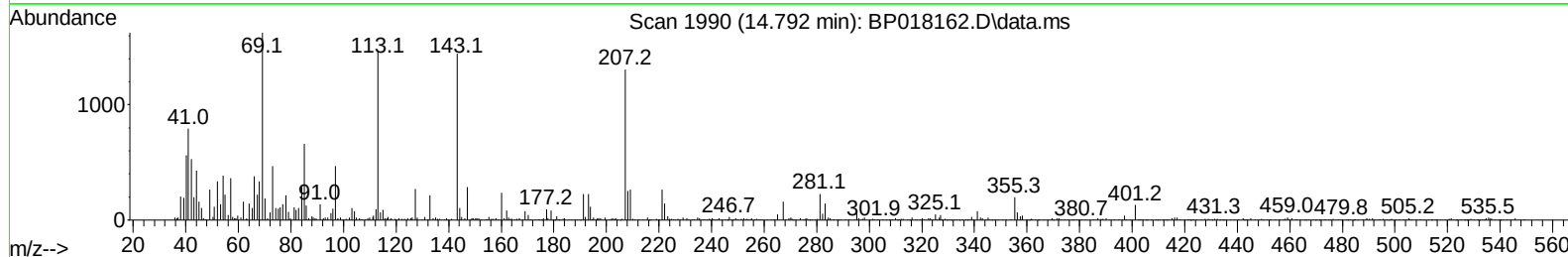
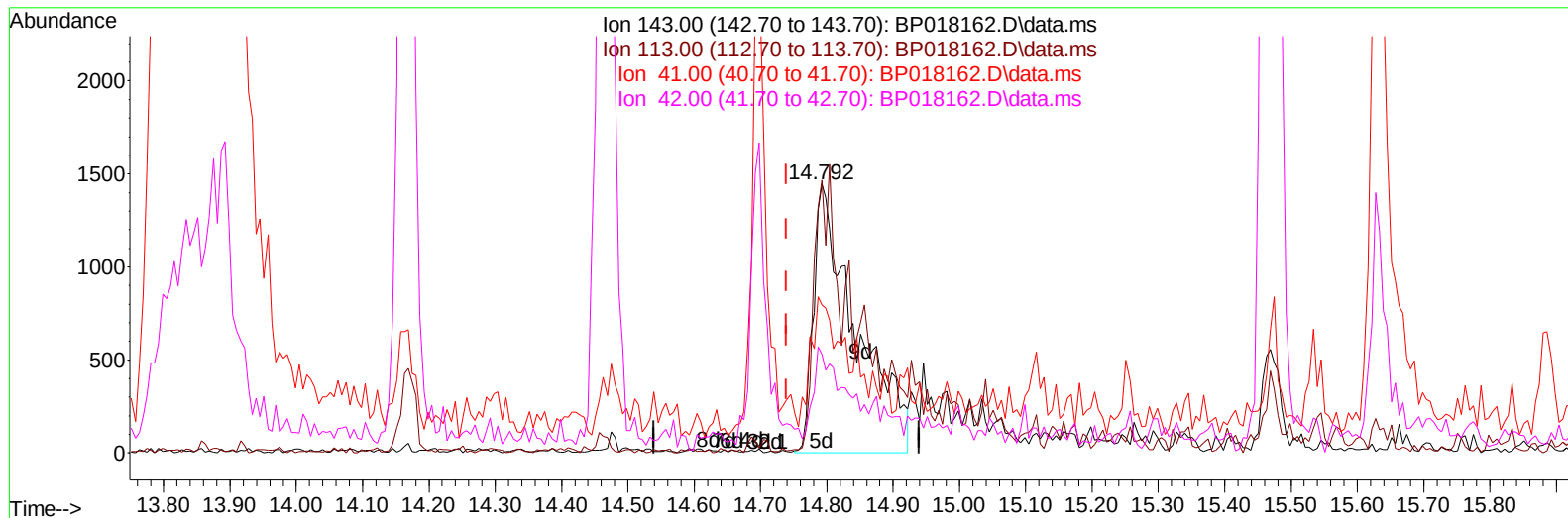
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(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.053) 2.31 ng/ul m

response 6448

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	101.59
41.00	47.70	55.02
42.00	30.10	36.75#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
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Quant Time: Nov 15 01:30:55 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	80536	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	326327	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	202337	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	453056	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	453460	20.000	ng/ul	-0.01
88) Perylene-d12	23.845	264	499720	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6899	3.022	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.969	99	31219	4.015	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	115116	25.194	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	102812	16.698	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	56025	8.793	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	67988	23.125	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	71525	21.474	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	113618	19.139	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	74251	9.121	ng/ul	0.01
46) Dimethylphthalate-d6	13.892	166	498864	26.721	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	483930	23.799	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	6448m	2.311	ng/ul	0.05
60) Fluorene-d10	15.469	176	406566	26.377	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	58336	17.811	ng/ul	-0.03
73) Anthracene-d10	17.351	188	679922	27.748	ng/ul	0.00
81) Pyrene-d10	19.604	212	873527	28.839	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.680	264	832012	28.634	ng/ul	-0.02

Target Compounds Qvalue

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

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