

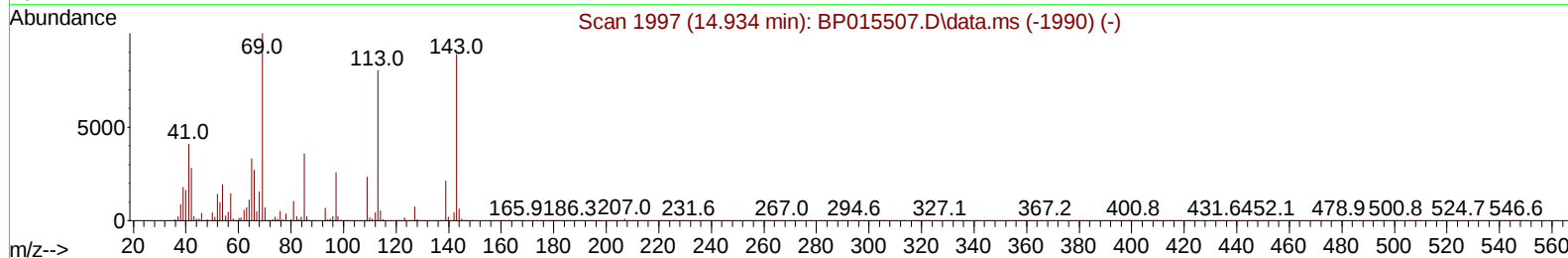
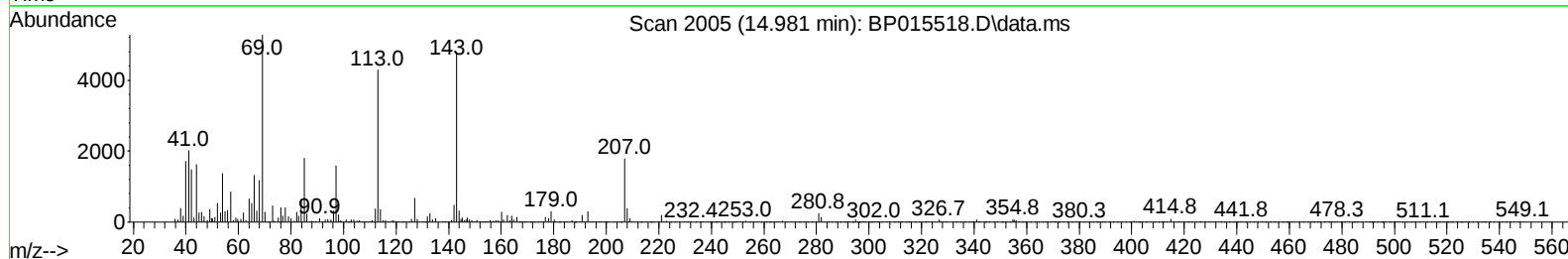
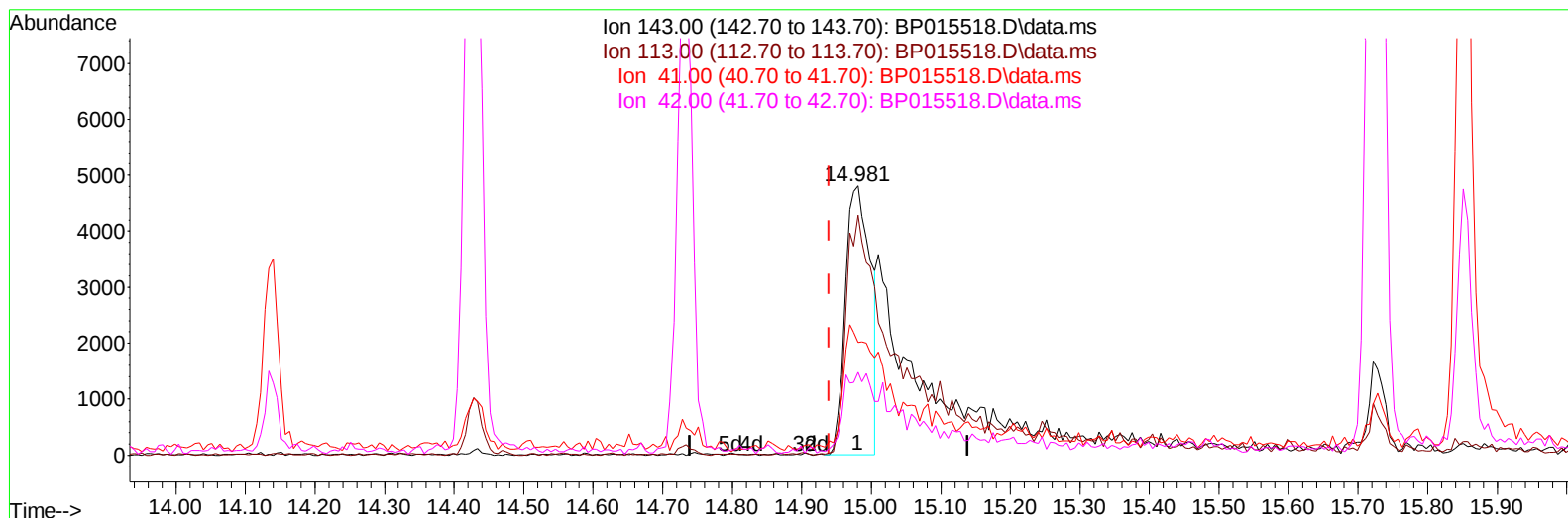
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015518.D
 Acq On : 14 Jun 2023 15:29
 Operator : MA/JU
 Sample : 03104-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONL8

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:09:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 09 23:33:38 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/15/2023
 Supervised By :mohammad ahmed 06/16/2023



TIC: BP015518.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.041) 2.14 ng/u1

response 12050

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	89.24
41.00	42.20	41.96
42.00	29.40	30.80

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Quant Time: Jun 14 23:40:32 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	115058	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	559177	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	397967	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.487	188	928407	20.000	ng/ul	0.00
79) Chrysene-d12	21.575	240	919039	20.000	ng/ul	0.00
88) Perylene-d12	24.122	264	965695	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	15203	4.893	ng/uL	0.00
4) Pyridine-d5	3.864	84	40688	5.041	ng/ul	0.02
7) Phenol-d5	7.246	99	83578	7.885	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	285174	45.048	ng/ul	0.00
11) 2-Chlorophenol-d4	7.617	132	264076	34.362	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	181942	20.914	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	190906	43.486	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	212299	42.346	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	325563	35.599	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	331708	25.269	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	1415092	45.092	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1548609	45.129	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143	21773m	3.874	ng/ul	0.04
60) Fluorene-d10	15.728	176	1198142	45.275	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	230566	39.202	ng/ul	0.00
73) Anthracene-d10	17.592	188	2038450	48.257	ng/ul	0.00
81) Pyrene-d10	19.816	212	2487793	50.577	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	2385762	49.237	ng/ul	0.00

Target Compounds Qvalue

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

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