

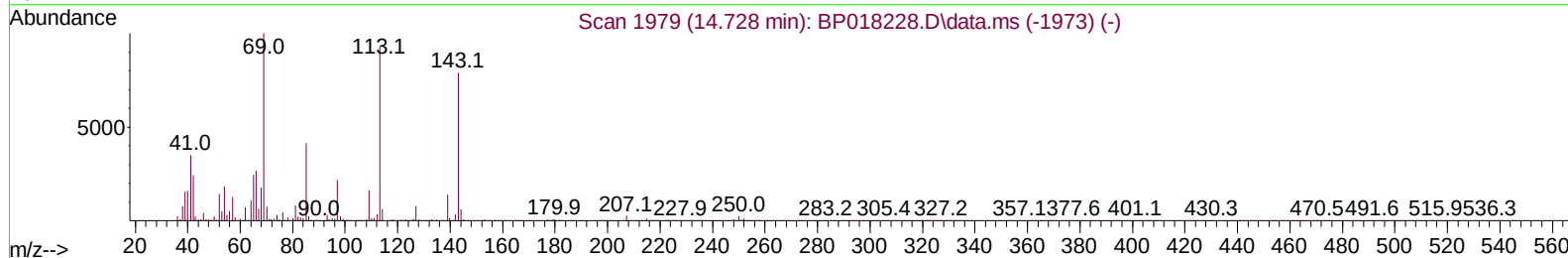
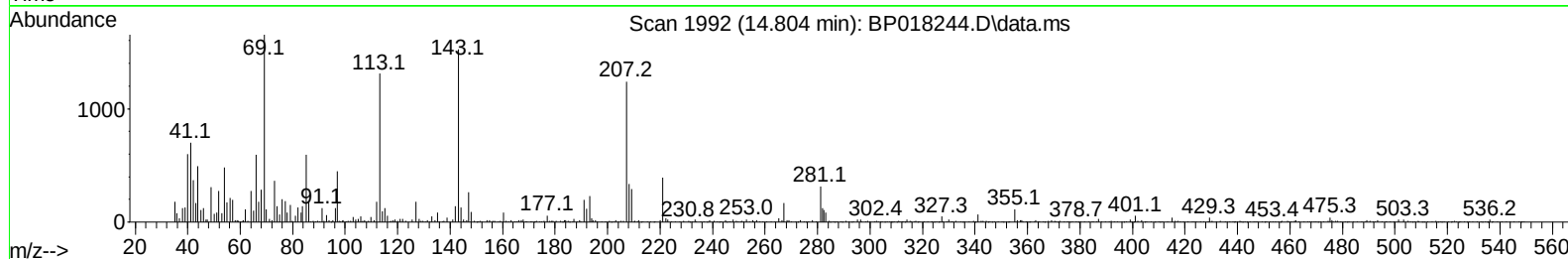
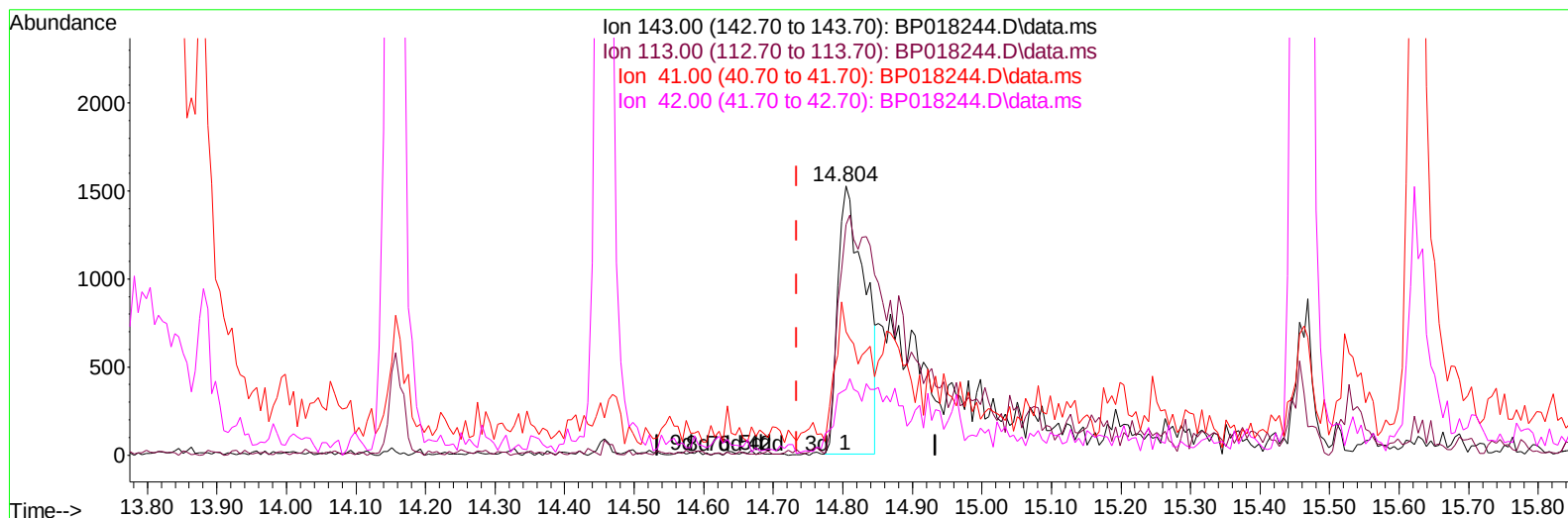
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018244.D
Acq On : 18 Nov 2023 04:46
Operator : MA/JU
Sample : 05369-15
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP2

Manual IntegrationsAPPROVED

Quant Time: Nov 19 00:45:24 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 18:13:37 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/19/2023
Supervised By :mohammad ahmed 11/19/2023



TIC: BP018244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.804min (+ 0.071) 1.41 ng/ul

response 4113

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	85.79
41.00	47.70	45.97
42.00	30.10	24.36

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Quant Time: Nov 19 01:03:12 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	73924	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	312841	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	211362	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	493700	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	485110	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	529416	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	9280	4.428	ng/uL	0.00
4) Pyridine-d5	3.599	84	59081	10.813	ng/ul	0.00
7) Phenol-d5	6.963	99	32561	4.562	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	123594	29.468	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	118439	20.957	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	74658	12.766	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	81061	28.760	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	88959	27.860	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	142518	25.042	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	198144	25.388	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	606699	31.110	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	632621	29.783	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	7428m	2.549	ng/ul	0.07
60) Fluorene-d10	15.463	176	511608	31.775	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	77149	21.616	ng/ul	0.00
73) Anthracene-d10	17.339	188	832356	31.173	ng/ul	0.00
81) Pyrene-d10	19.598	212	1028218	31.732	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	991791	32.218	ng/ul	0.00

Target Compounds Qvalue

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

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