

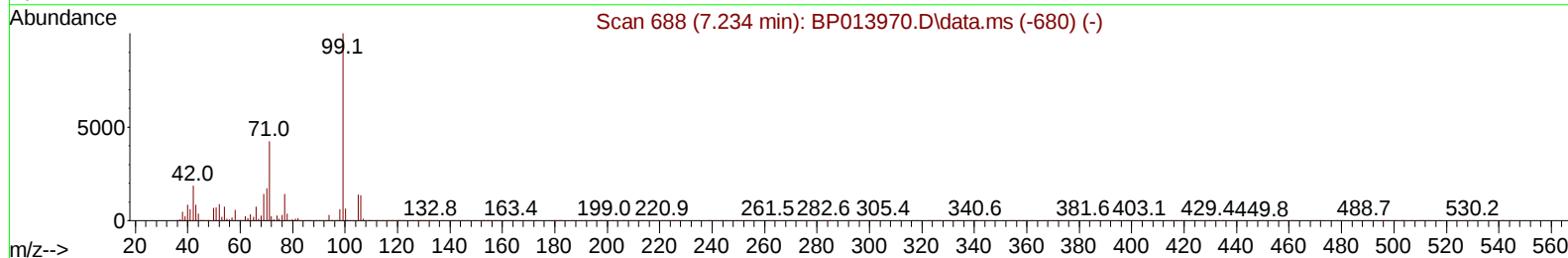
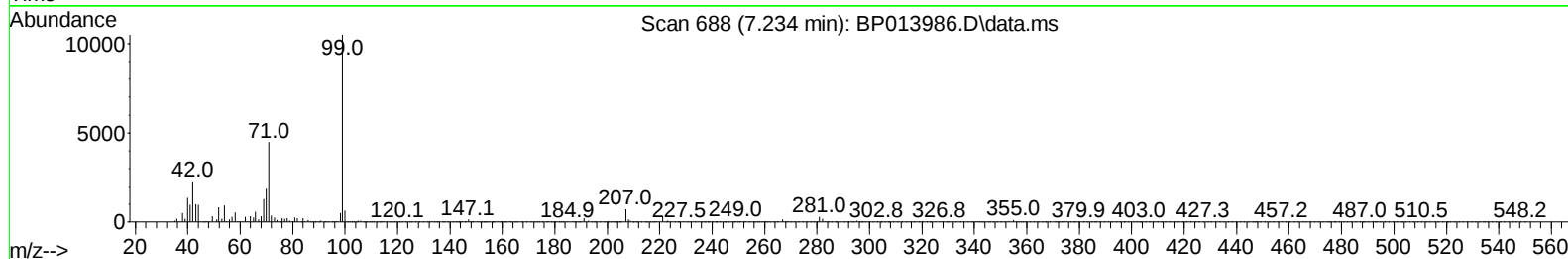
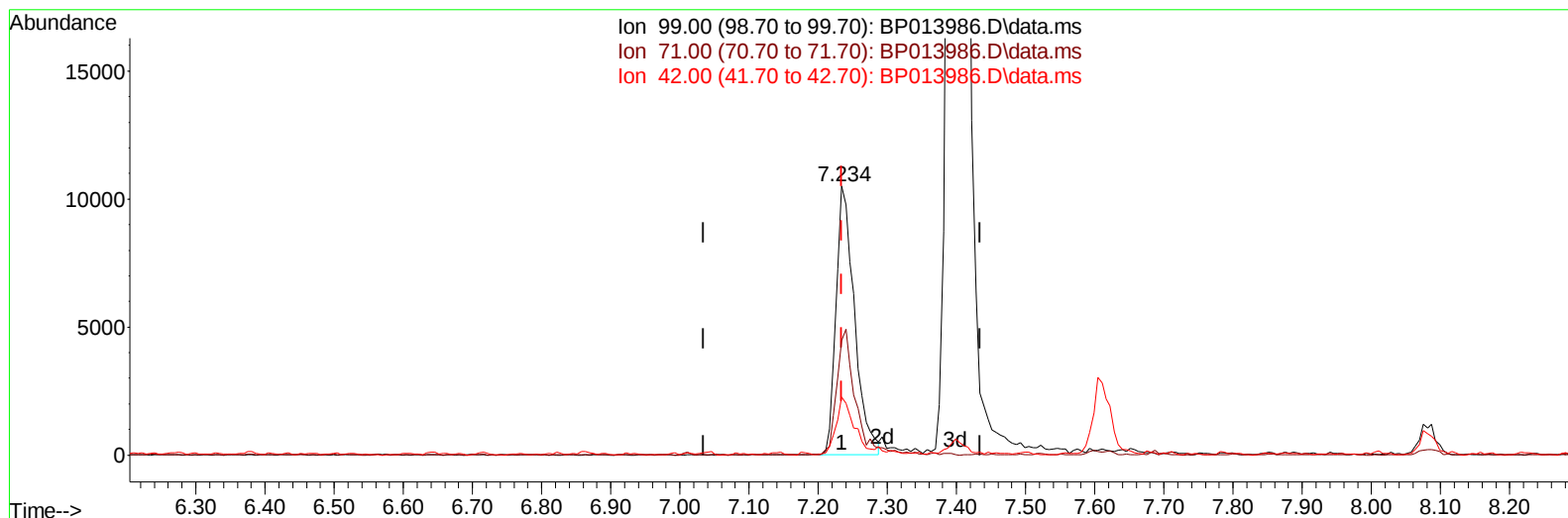
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013986.D
Acq On : 08 Mar 2023 20:02
Operator : CG/JU
Sample : 01729-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BY5

Manual IntegrationsAPPROVED

Quant Time: Mar 09 01:17:02 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 09 01:10:12 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/09/2023
Supervised By :Jagrut Upadhyay 03/09/2023



TIC: BP013986.D\data.ms

(7) Phenol-d5 (S)

7.234min (-0.000) 2.30 ng/u1

response 19432

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	42.63
42.00	19.50	21.56
0.00	0.00	0.00

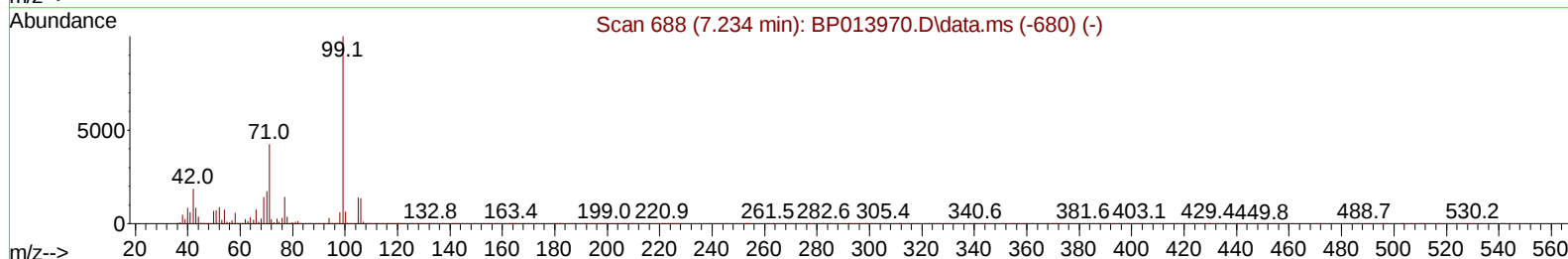
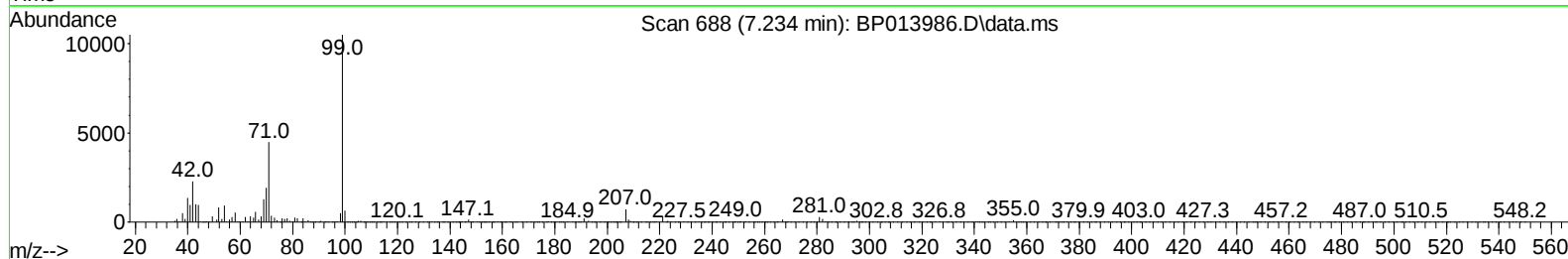
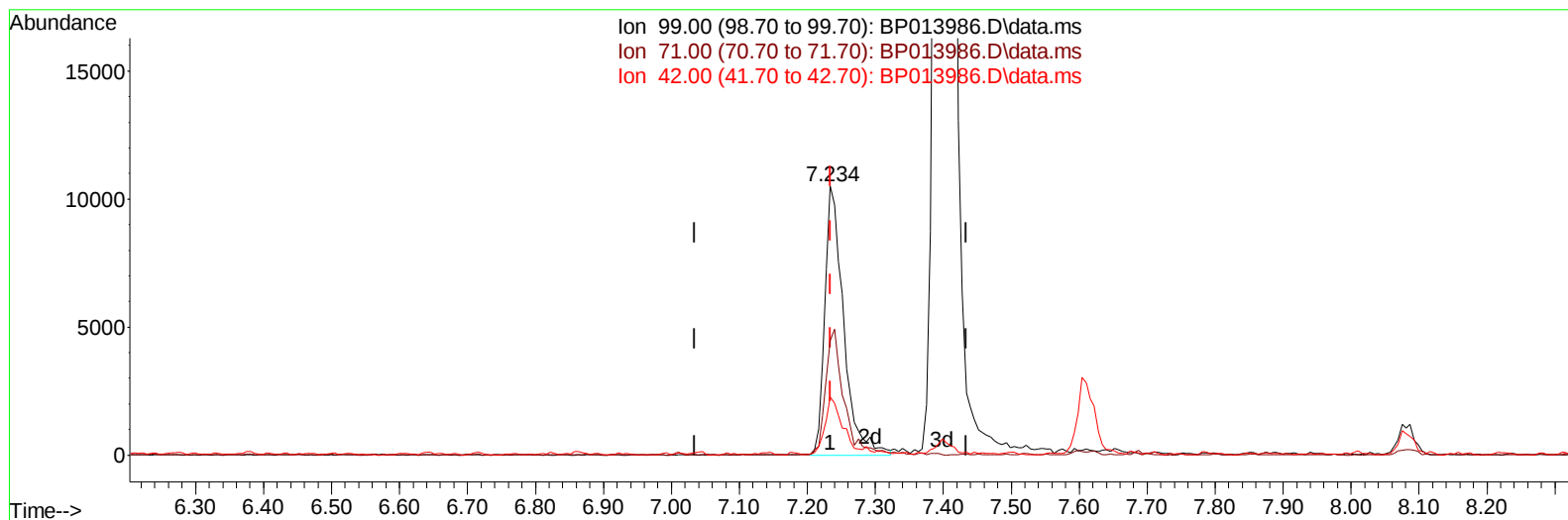
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TIC: BP013986.D\data.ms

(7) Phenol-d5 (S)

7.234min (-0.000) 2.39 ng/ul m

response 20219

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	42.63
42.00	19.50	21.56
0.00	0.00	0.00

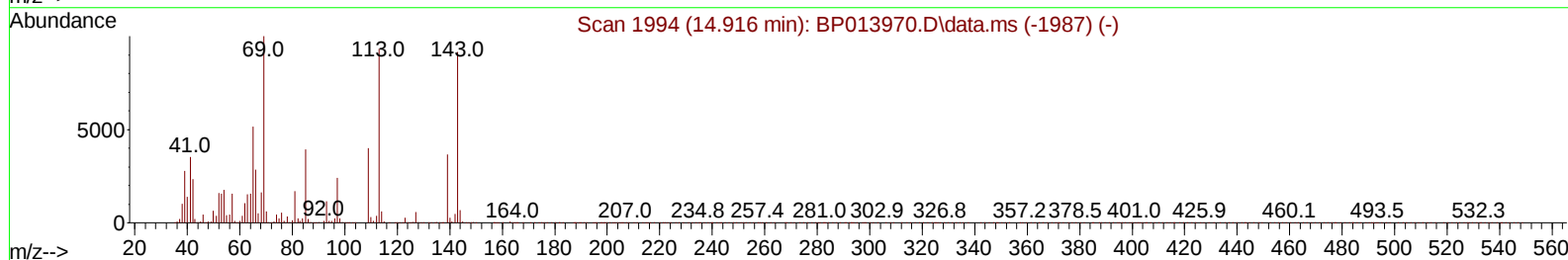
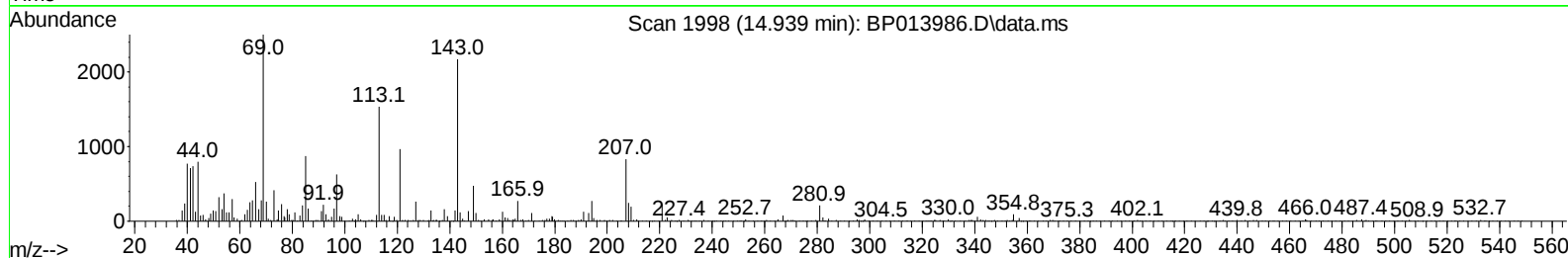
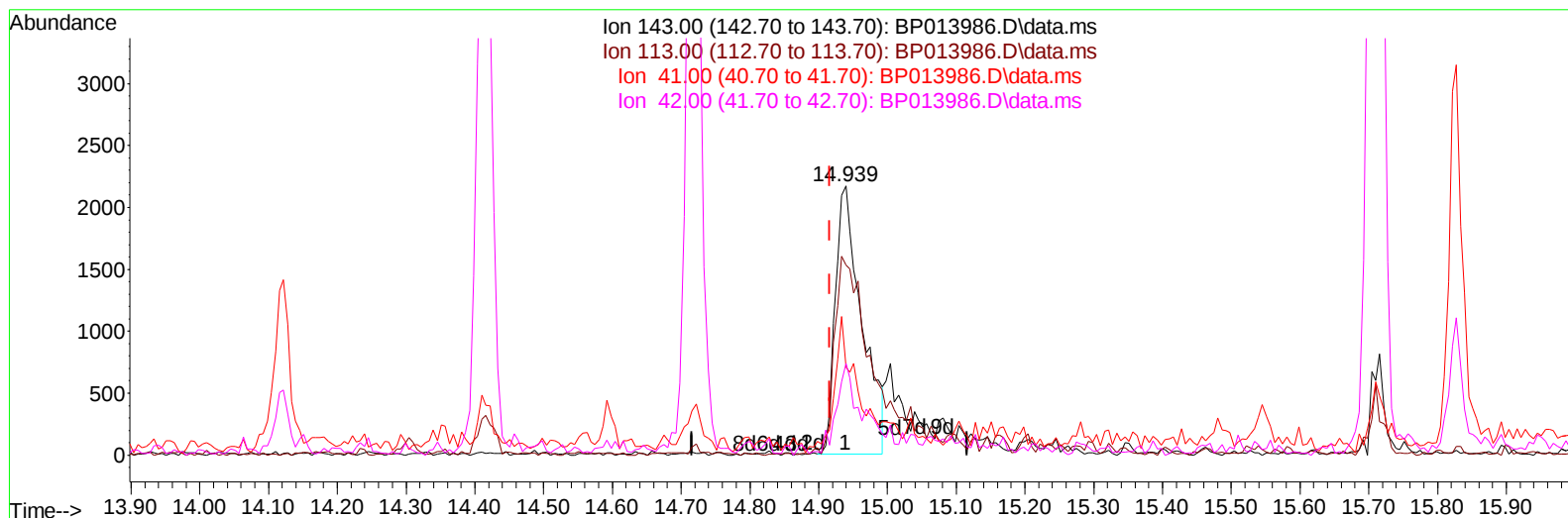
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
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 Acq On : 08 Mar 2023 20:02
 Operator : CG/JU
 Sample : 01729-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BY5

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TIC: BP013986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (+ 0.023) 1.44 ng/ul

response 5766

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	70.63#
41.00	40.20	32.78
42.00	28.50	34.12

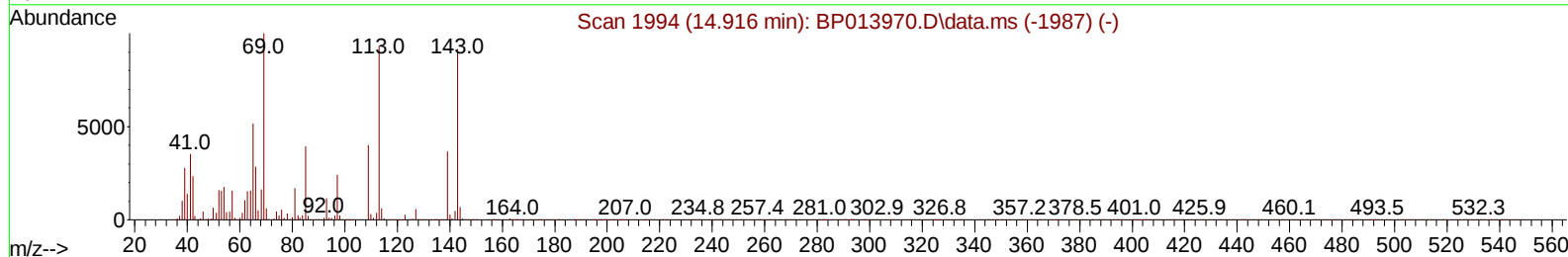
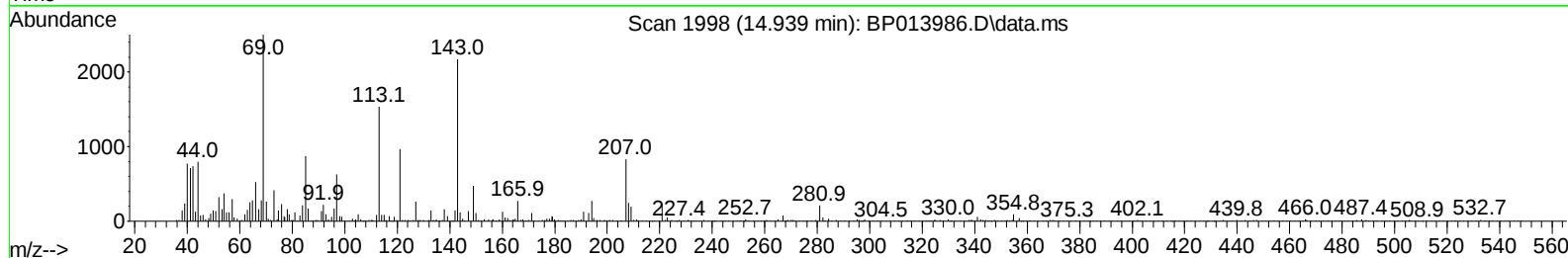
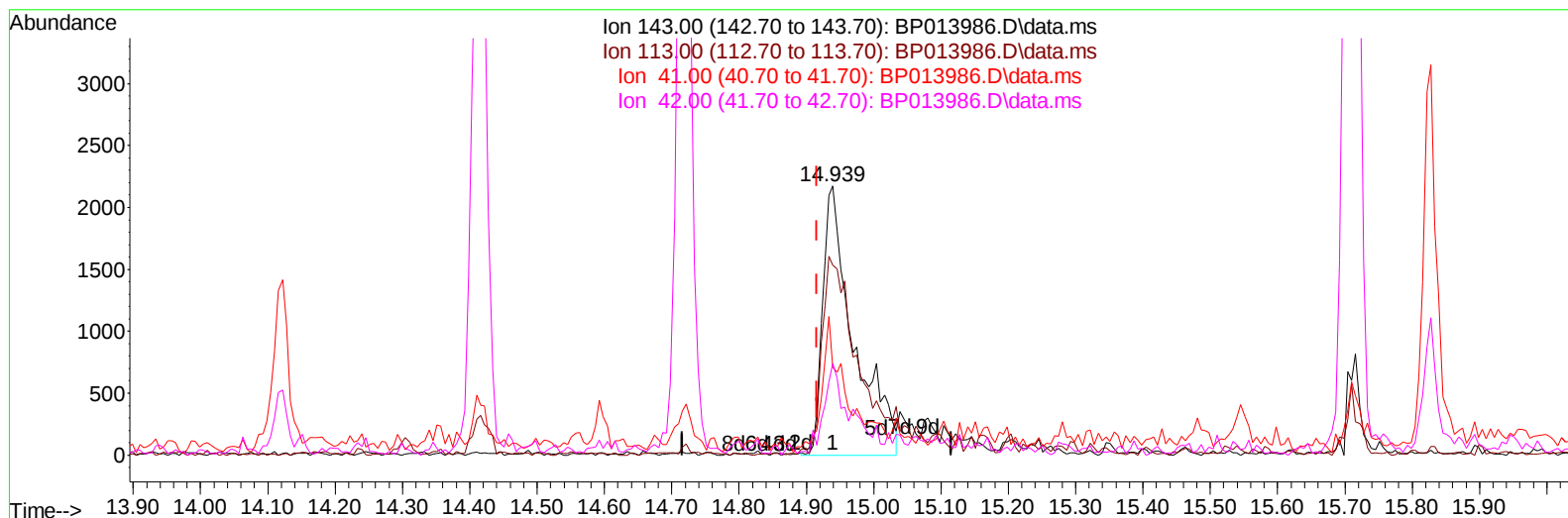
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013986.D
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Sample : 01729-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP013986.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (+ 0.023) 1.72 ng/ul m

response 6917

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	70.63#
41.00	40.20	32.78
42.00	28.50	34.12

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 ALS Vial : 17 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	95682	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	375770	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.722	164	268297	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.475	188	689585	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	839466	20.000	ng/ul	0.00
88) Perylene-d12	24.086	264	967264	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	10501	4.147	ng/uL	0.00
4) Pyridine-d5	3.875	84	54827	7.819	ng/ul	0.00
7) Phenol-d5	7.234	99	20219m	2.395	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	111251	22.592	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	67438	10.483	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	48090	6.972	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128	68868	22.969	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	46923	13.074	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	87142	12.904	ng/ul	0.00
31) 4-Chloroaniline-d4	11.046	131	121317	13.575	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	593041	26.045	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	589242	24.429	ng/ul	0.00
54) 4-Nitrophenol-d4	14.939	143	6917m	1.724	ng/ul	0.02
60) Fluorene-d10	15.710	176	515656	26.722	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	61815	11.860	ng/ul	0.00
73) Anthracene-d10	17.575	188	989905	29.571	ng/ul	0.00
81) Pyrene-d10	19.792	212	1322092	28.555	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	1525905	29.657	ng/ul	0.00

Target Compounds Qvalue

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

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