

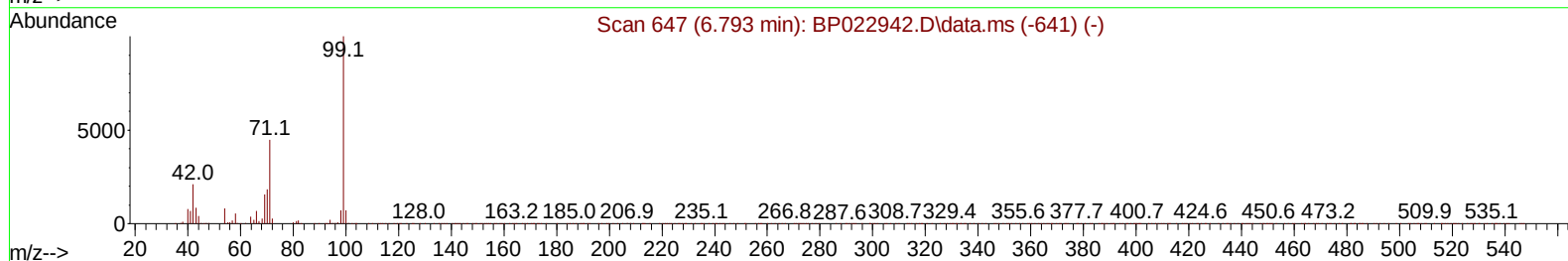
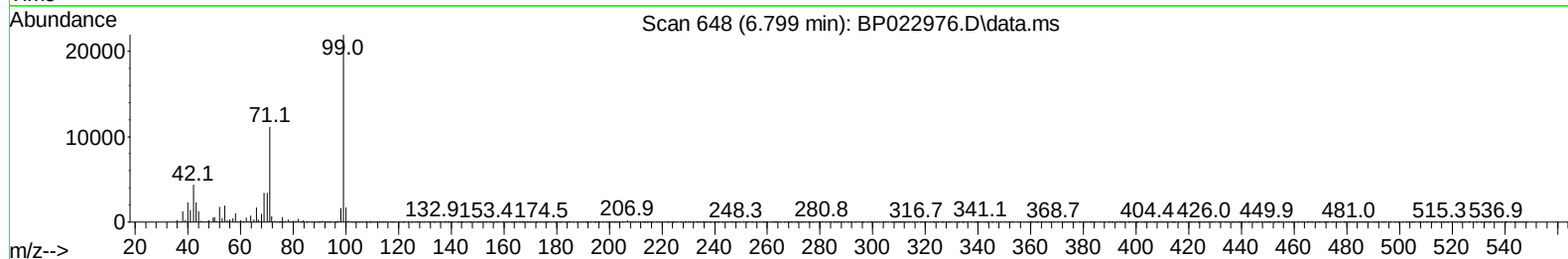
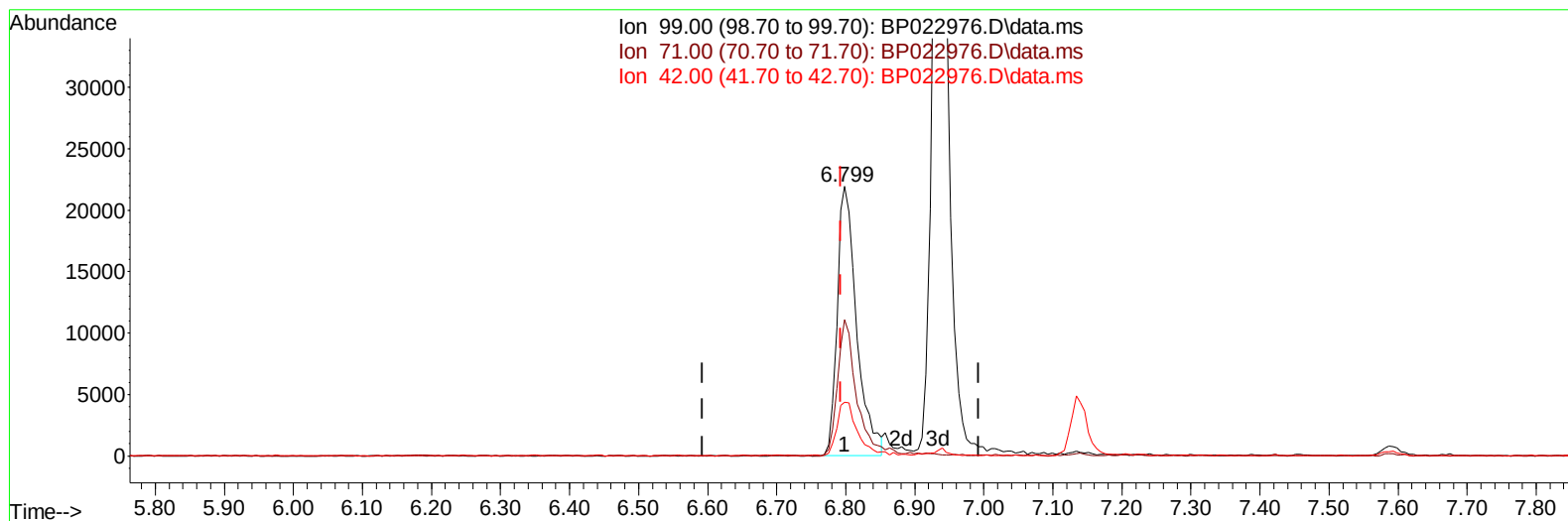
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022976.D
Acq On : 09 Nov 2024 21:18
Operator : RC/JU
Sample : P4746-21
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP94

Manual IntegrationsAPPROVED

Quant Time: Nov 09 23:35:06 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 08 23:49:48 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/11/2024
Supervised By :mohammad ahmed 11/11/2024



TIC: BP022976.D\data.ms

(7) Phenol-d5 (S)

6.799min (+ 0.006) 9.66 ng/u1

response 42795

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	50.70
42.00	21.10	19.98
0.00	0.00	0.00

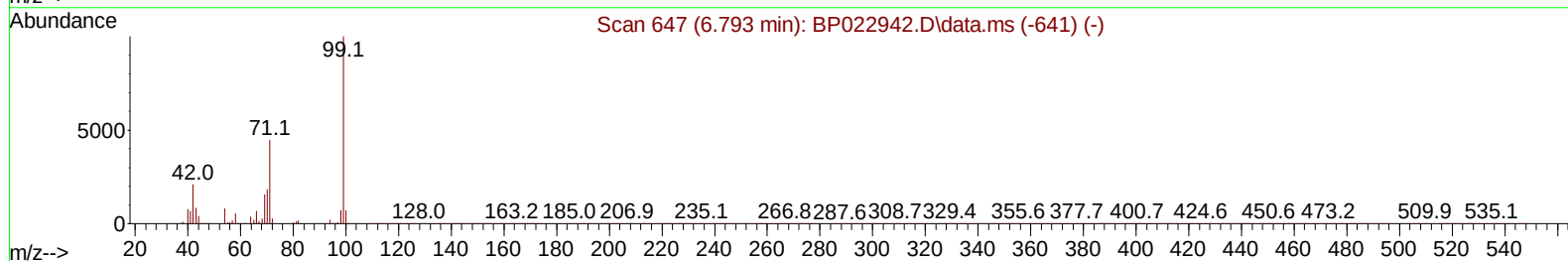
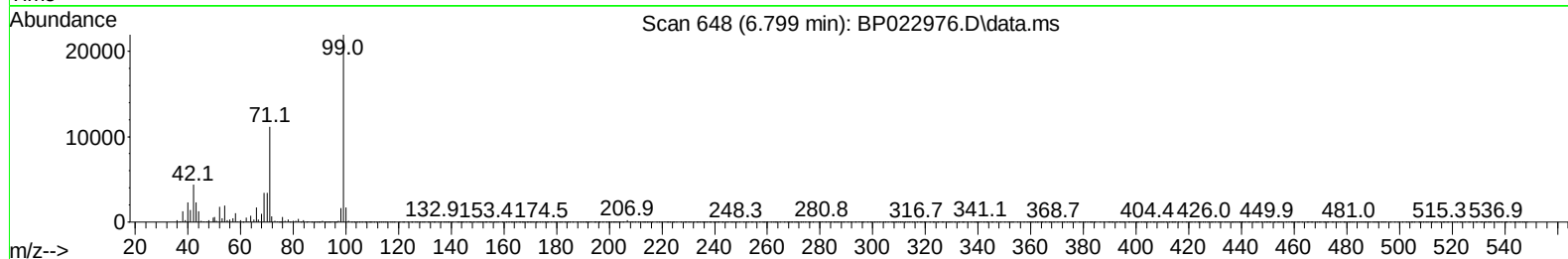
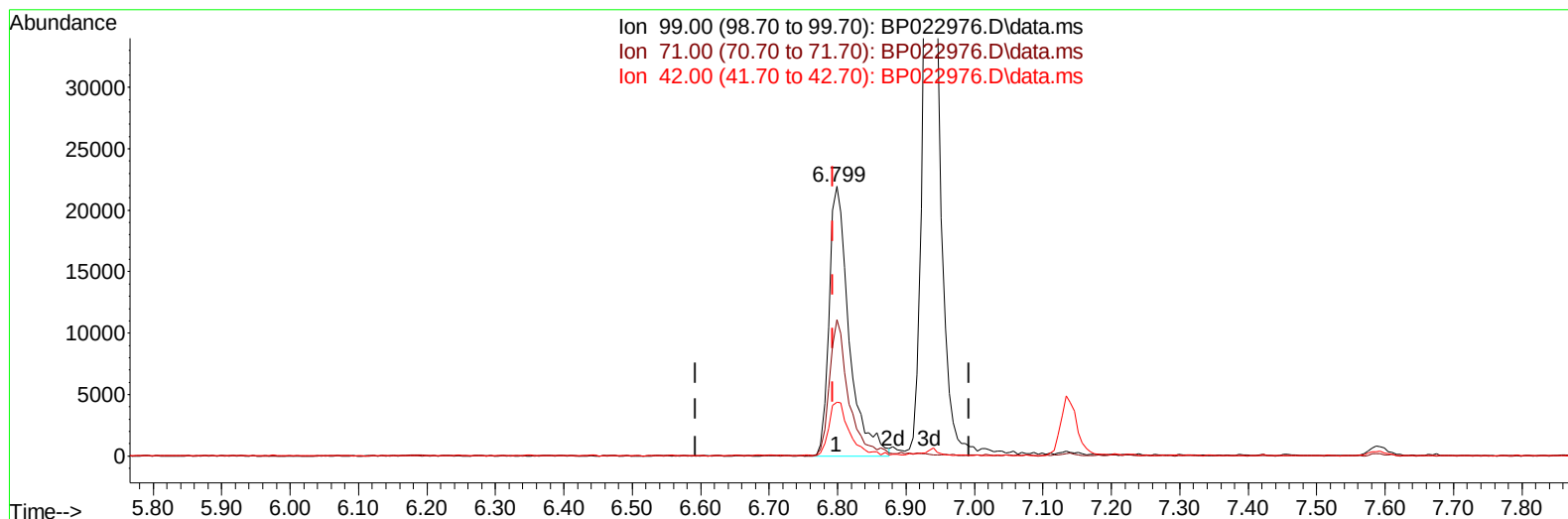
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(7) Phenol-d5 (S)

6.799min (+ 0.006) 10.00 ng/ul m

response 44319

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	50.70
42.00	21.10	19.98
0.00	0.00	0.00

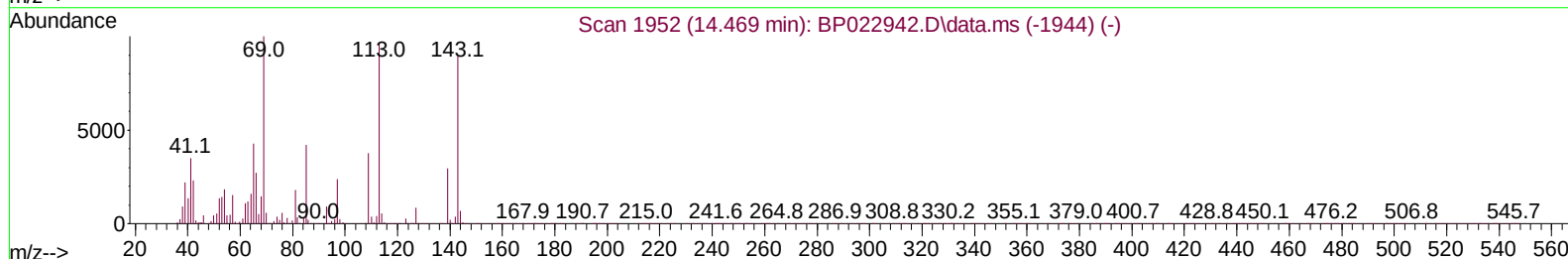
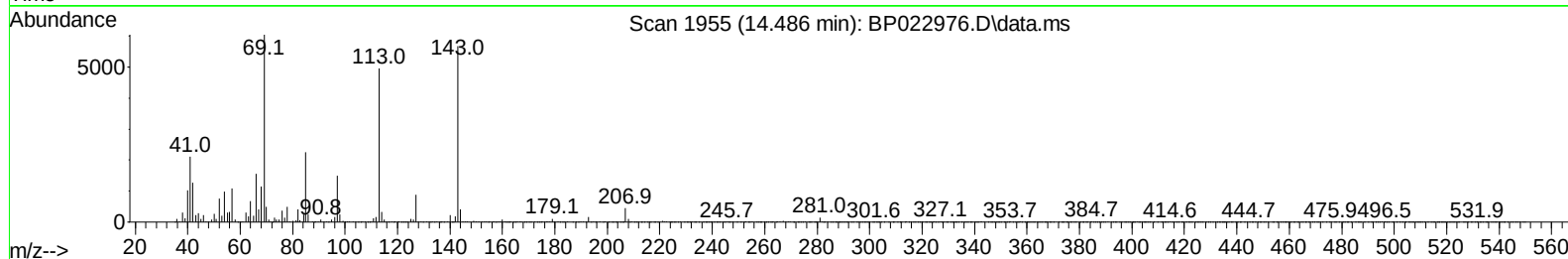
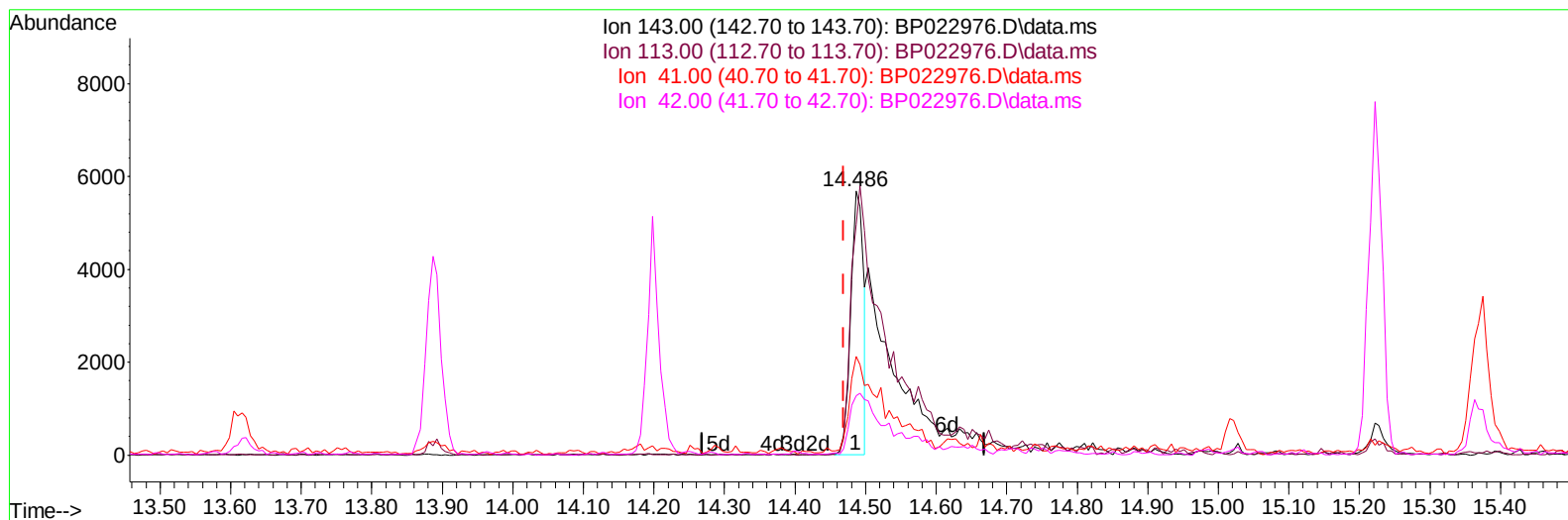
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TIC: BP022976.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.018) 3.64 ng/ul

response 7196

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	87.44
41.00	39.20	37.28
42.00	25.90	22.40

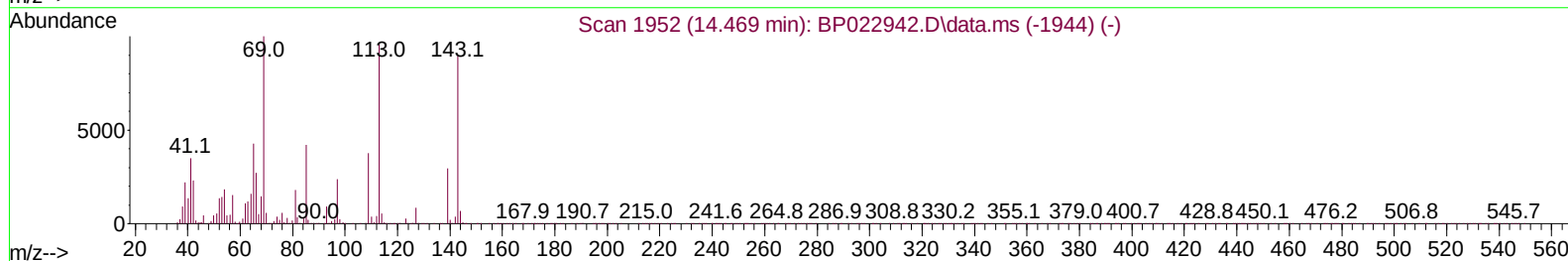
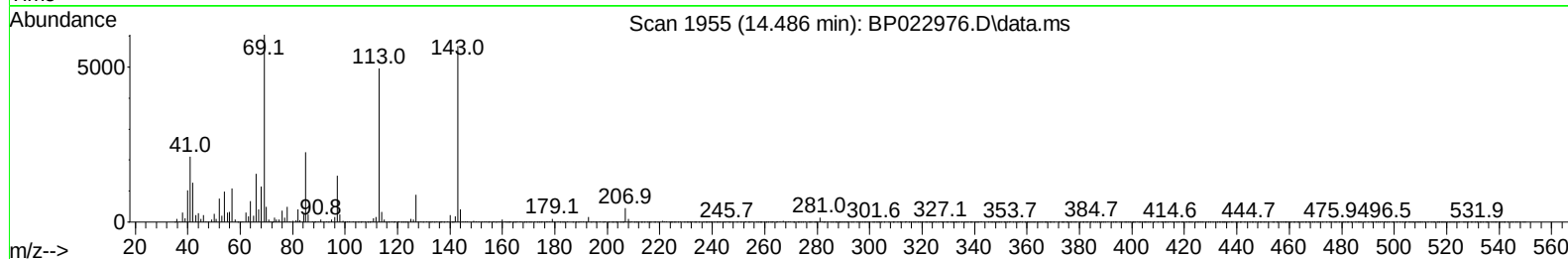
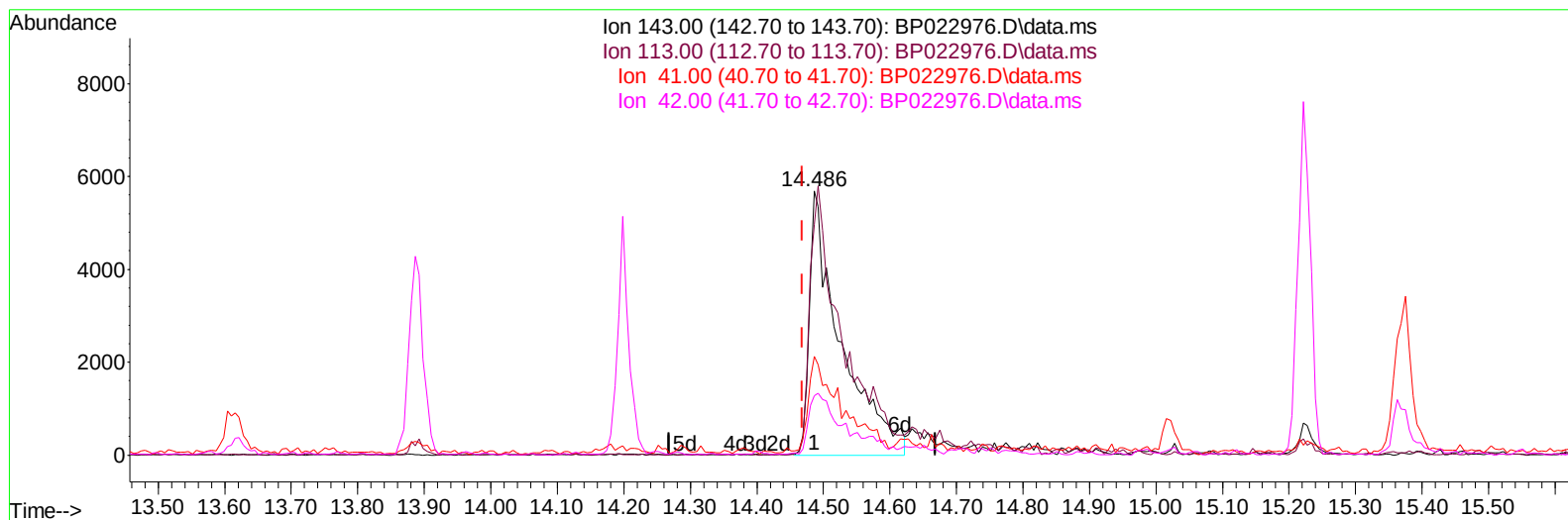
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 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
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TIC: BP022976.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.018) 9.36 ng/ul m

response 18522

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	87.44
41.00	39.20	37.28
42.00	25.90	22.40

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/11/2024
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Quant Time: Nov 10 00:27:54 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	59929	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.340	136	227144	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.198	164	182680	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.016	188	474295	20.000	ng/ul	-0.02
79) Chrysene-d12	21.451	240	572071	20.000	ng/ul	0.00
88) Perylene-d12	24.686	264	700849	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	6112	4.902	ng/uL	0.00
4) Pyridine-d5	3.563	84	29236	7.893	ng/ul	0.00
7) Phenol-d5	6.799	99	44319m	10.003	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	80447	30.904	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.140	132	94541	29.135	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	78078	21.414	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	51416	32.515	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	62991	33.413	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	122594	31.052	ng/ul	0.00
31) 4-Chloroaniline-d4	10.487	131	126931	26.950	ng/ul	-0.01
46) Dimethylphthalate-d6	13.616	166	471311	35.797	ng/ul	0.00
49) Acenaphthylene-d8	13.886	160	460277	33.504	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.486	143	18522m	9.361	ng/ul	0.02
60) Fluorene-d10	15.222	176	411645	35.660	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.369	200	88123	31.015	ng/ul	0.00
73) Anthracene-d10	17.110	188	718434	35.775	ng/ul	-0.01
81) Pyrene-d10	19.504	212	985079	37.384	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.457	264	1273068	38.134	ng/ul	-0.01

Target Compounds Qvalue

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

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