

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
Data File : BG063304.D
Acq On : 12 Nov 2024 00:43
Operator : RC/JU
Sample : P4644-17
Misc :
ALS Vial : 22 Sample Multiplier: 1

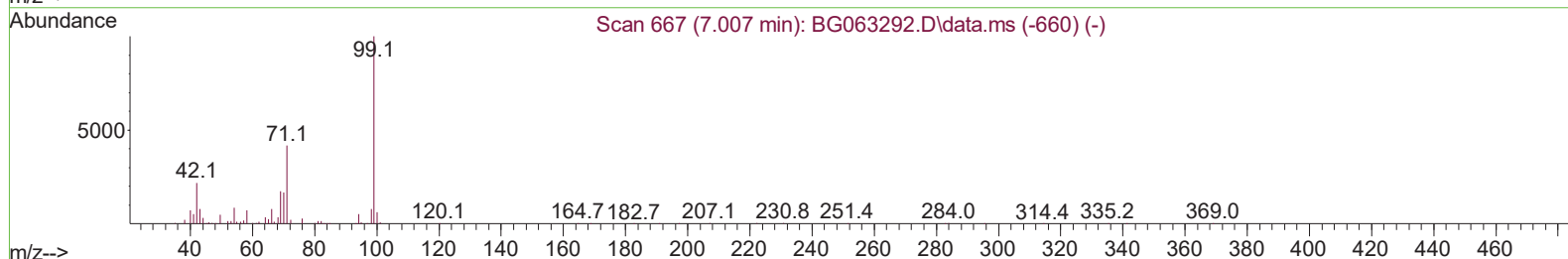
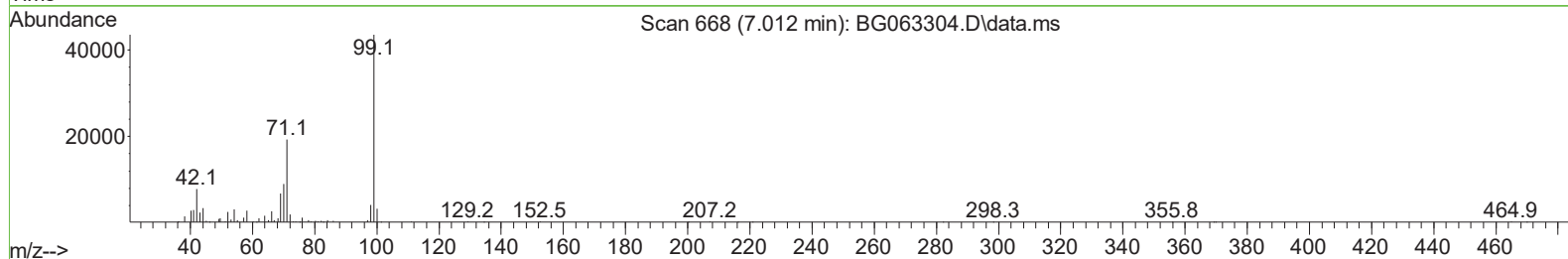
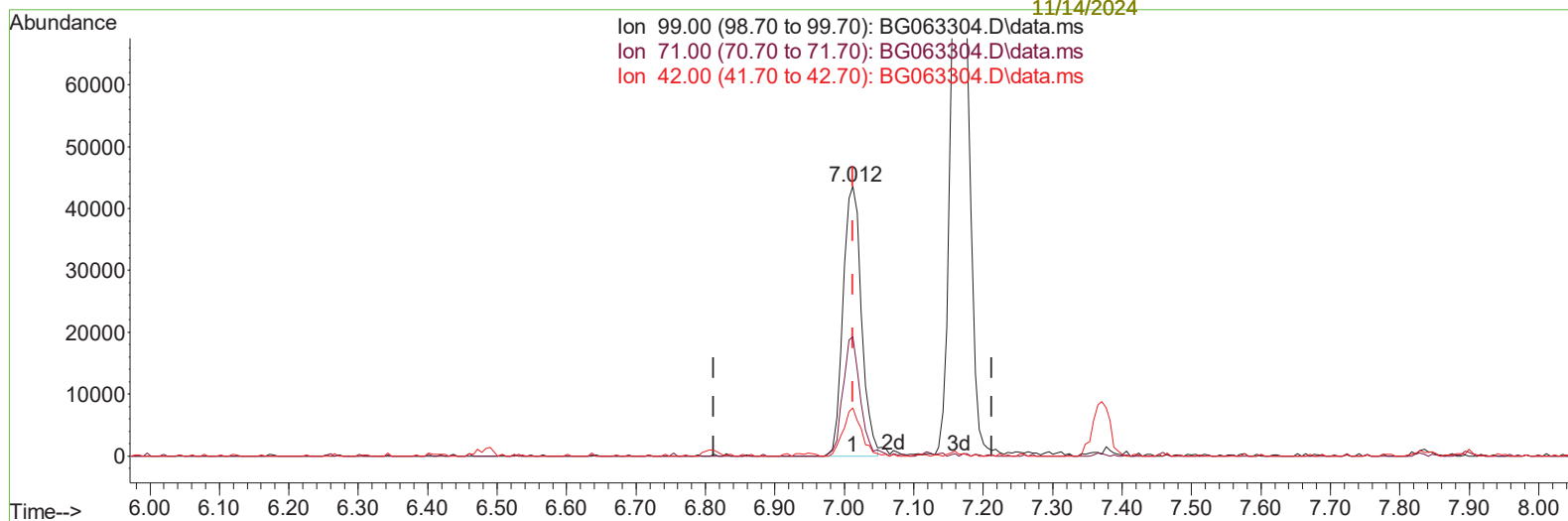
Instrument :
BNA_G
ClientSampleId :
C0CH2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/12/2024
Supervised By :mohammad ahmed 11/14/2024

Supervised By :mohammad
ahmed
11/14/2024

Quant Time: Nov 12 00:37:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration



TIC: BG063304.D\data.ms

(7) Phenol-d5 (S)

7.012min (-0.000) 8.69 ng/u1

response 79739

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	48.00	44.17
42.00	19.70	17.84
0.00	0.00	0.00

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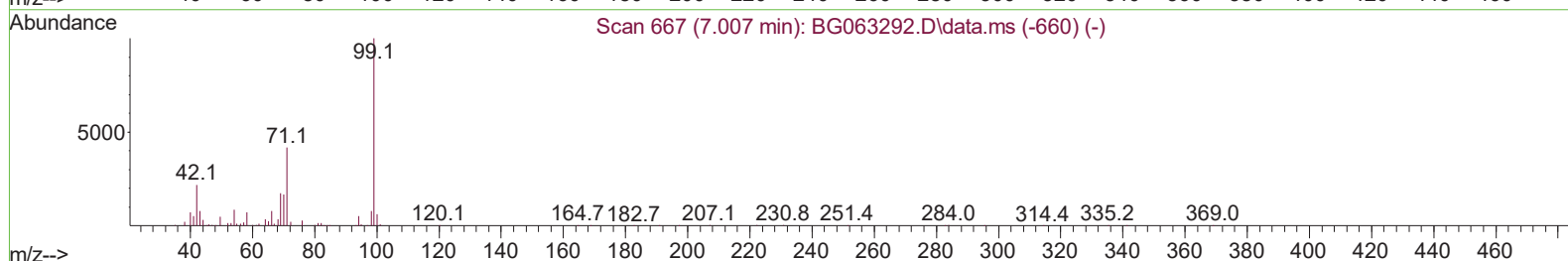
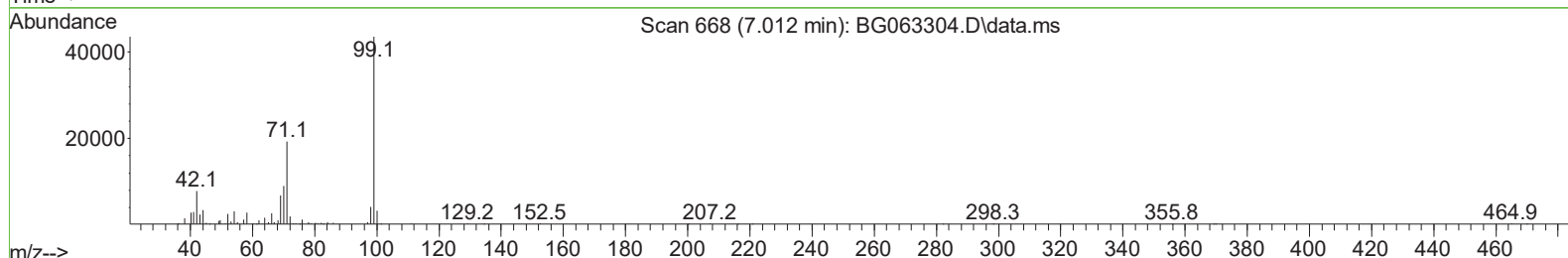
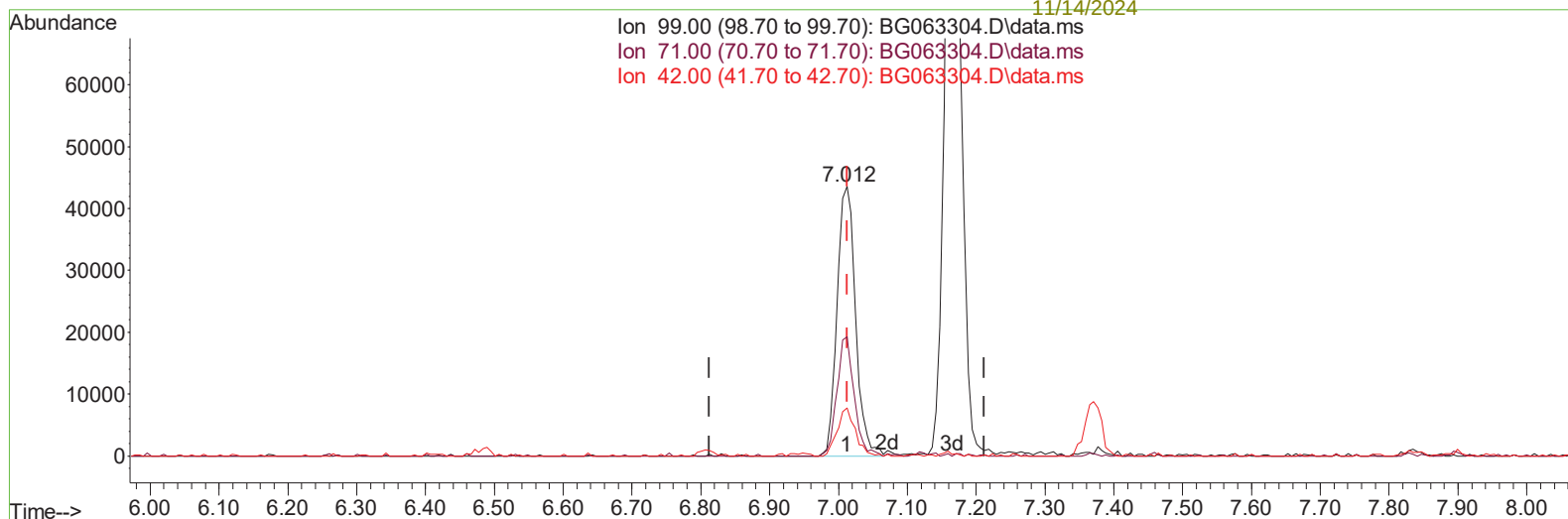
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(7) Phenol-d5 (S)

7.012min (-0.000) 8.80 ng/ul m

response 80741

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	48.00	44.17
42.00	19.70	17.84
0.00	0.00	0.00

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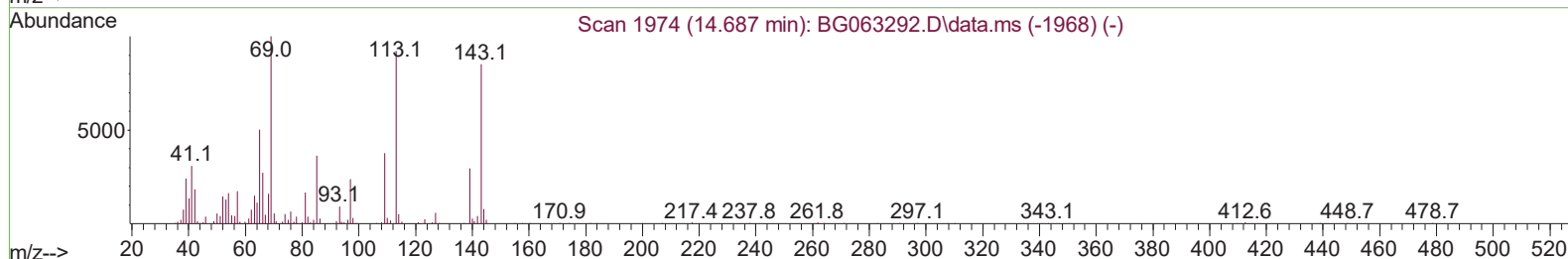
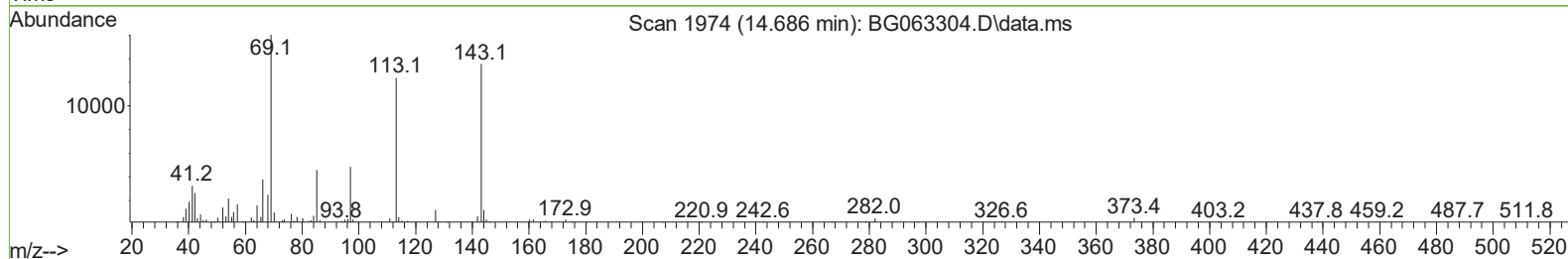
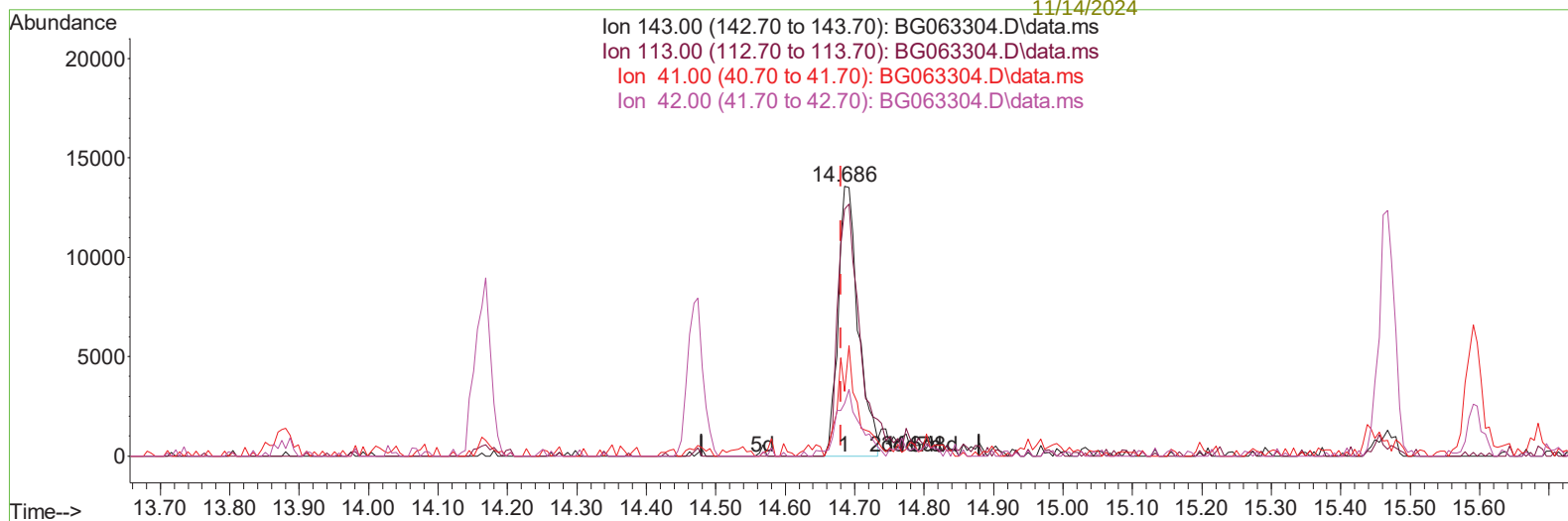
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(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 6.12 ng/u1

response 27452

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	91.51#
41.00	42.60	23.95#
42.00	30.90	19.39#

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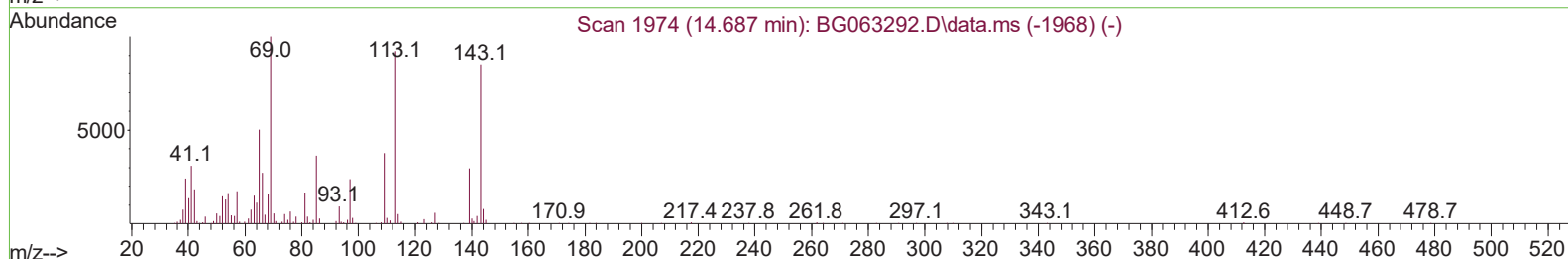
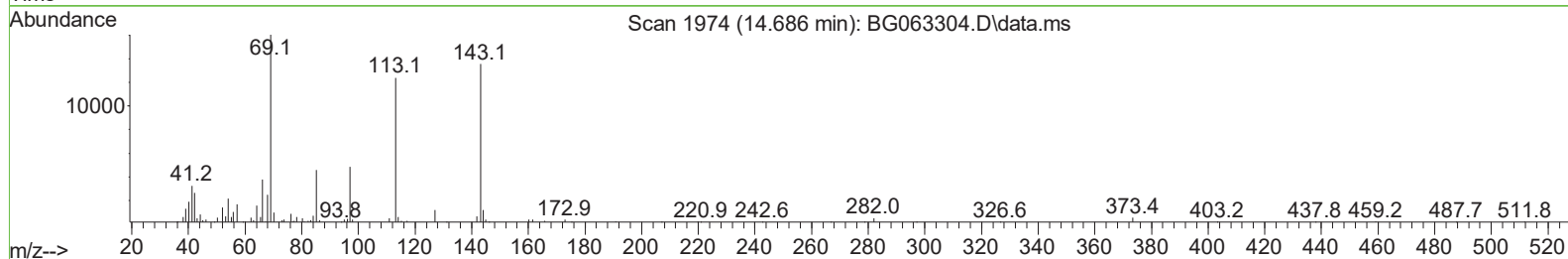
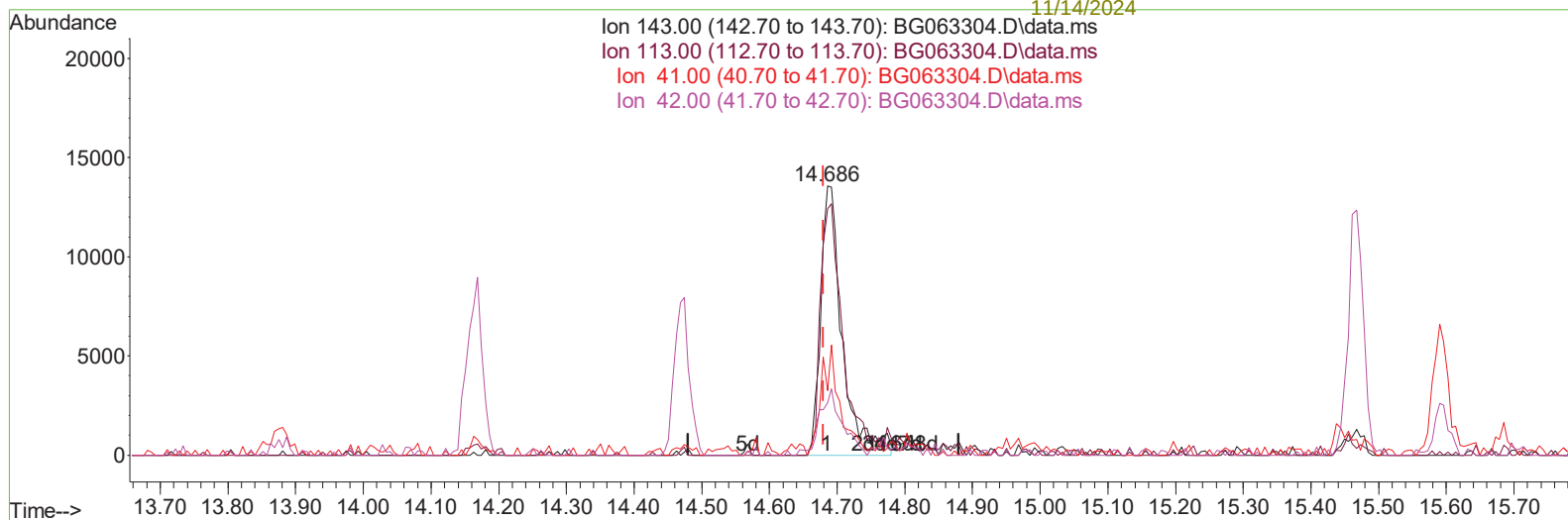
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(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 6.70 ng/ul m

response 30089

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	91.51#
41.00	42.60	23.95#
42.00	30.90	19.39#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.835	152	101728	20.000	ng/ul	0.00
20) Naphthalene-d8	10.626	136	488248	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.474	164	418562	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	984332	20.000	ng/ul	0.00
79) Chrysene-d12	21.478	240	1005592	20.000	ng/ul	0.00
88) Perylene-d12	24.515	264	1078619	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	11242	4.997	ng/uL	0.00
4) Pyridine-d5	3.716	84	50973	6.855	ng/ul	0.00
7) Phenol-d5	7.012	99	80741m	8.796	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	190290	35.342	ng/ul	0.00
11) 2-Chlorophenol-d4	7.371	132	189350	31.780	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	170313	21.890	ng/ul	0.00
21) Nitrobenzene-d5	8.986	128	125057	36.433	ng/ul	0.00
24) 2-Nitrophenol-d4	9.709	143	156371	35.248	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.250	165	302708	34.769	ng/ul	0.00
31) 4-Chloroaniline-d4	10.761	131	263608	23.246	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1062250	32.399	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1128885	35.178	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	30089m	6.703	ng/ul	0.00
60) Fluorene-d10	15.467	176	957887	34.618	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.596	200	199803	32.298	ng/ul	0.00
73) Anthracene-d10	17.324	188	1597540	37.762	ng/ul	0.00
81) Pyrene-d10	19.621	212	2045097	40.563	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.310	264	2043147	39.233	ng/ul	0.00

Target Compounds Qvalue

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

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