

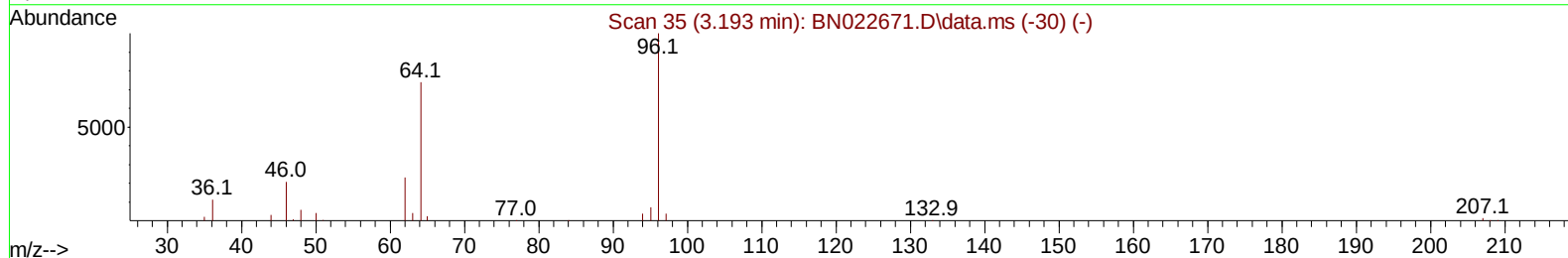
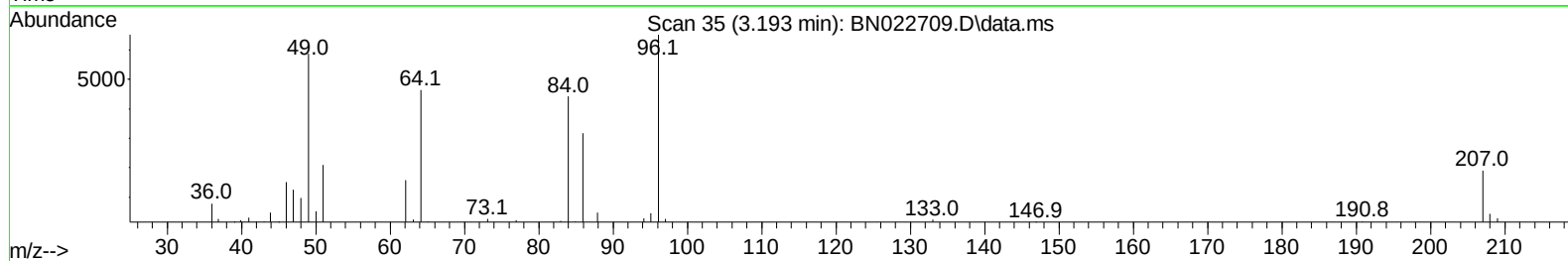
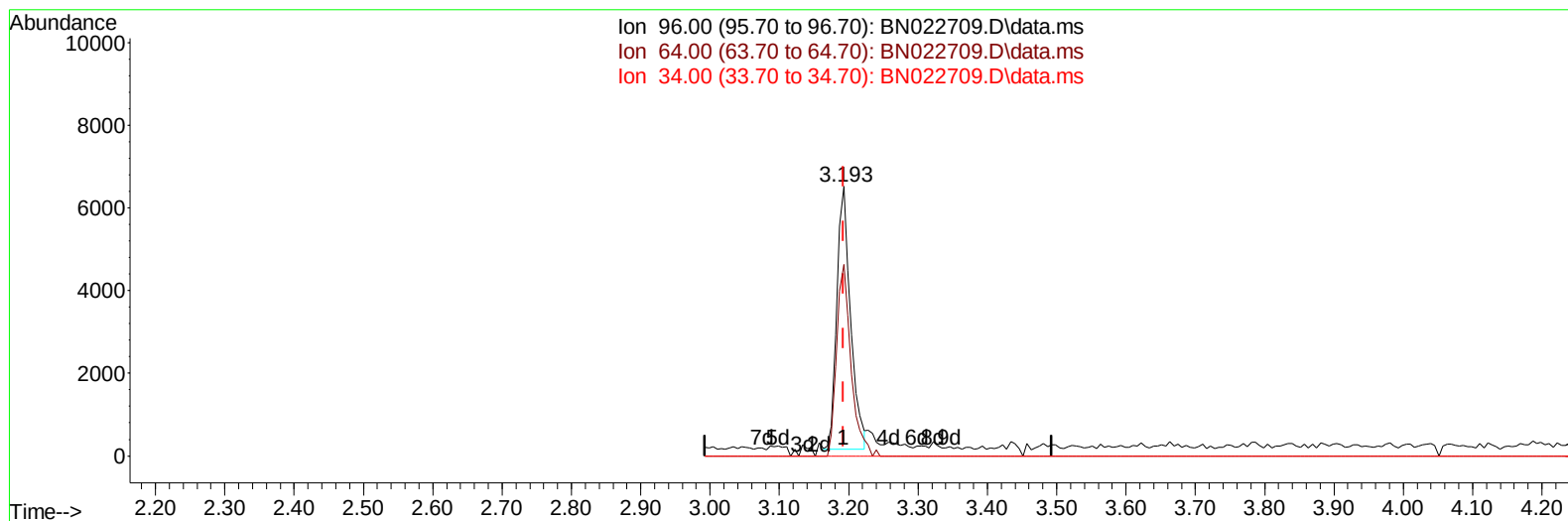
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111522\
 Data File : BN022709.D
 Acq On : 17 Nov 2022 05:12
 Operator : CG/JU
 Sample : N5577-06
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C9819

Manual IntegrationsAPPROVED

Quant Time: Nov 17 05:56:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 15 22:53:43 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/17/2022
 Supervised By :mohammad ahmed 11/18/2022



TIC: BN022709.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.193min (-0.000) 3.69 ng/uL

response 8772

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	71.15
34.00	0.00	0.00
0.00	0.00	0.00

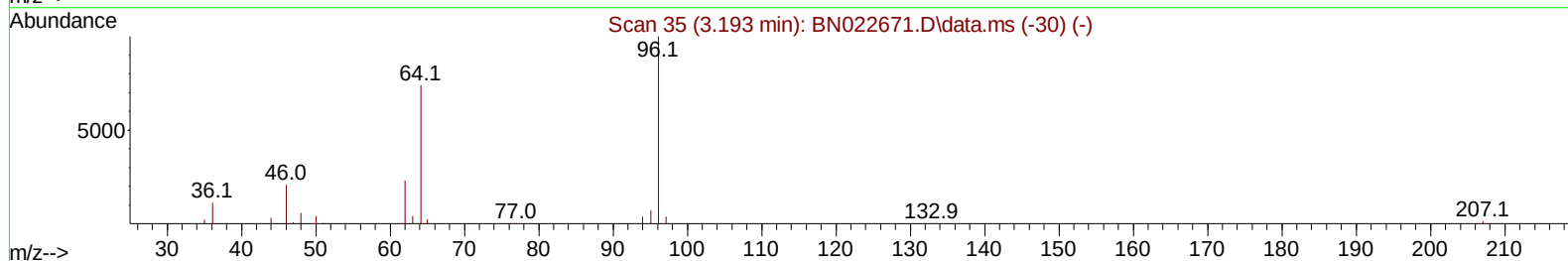
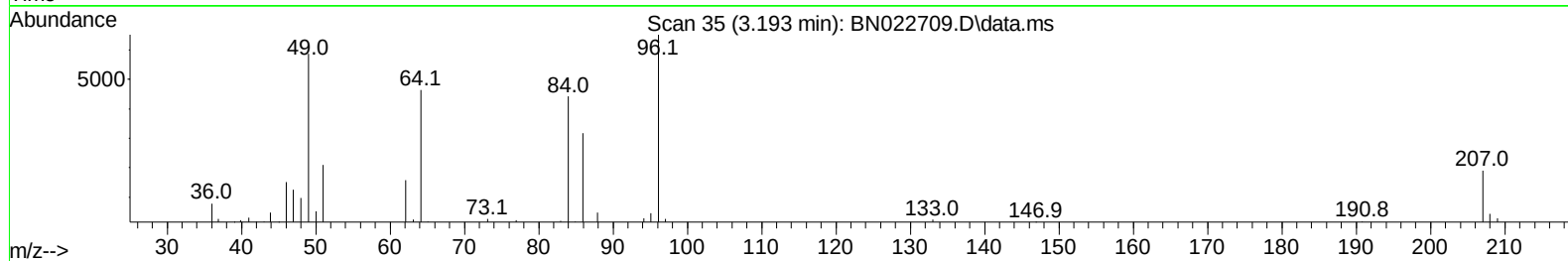
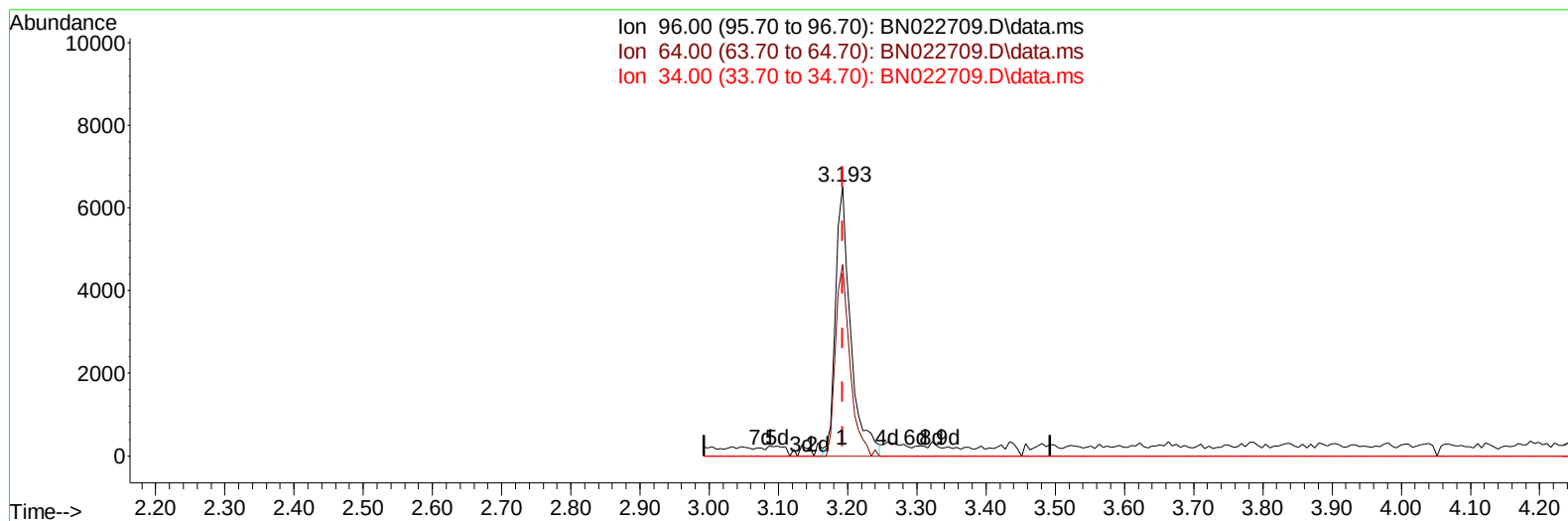
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TIC: BN022709.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.193min (-0.000) 4.20 ng/uL m

response 9969

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	71.15
34.00	0.00	0.00
0.00	0.00	0.00

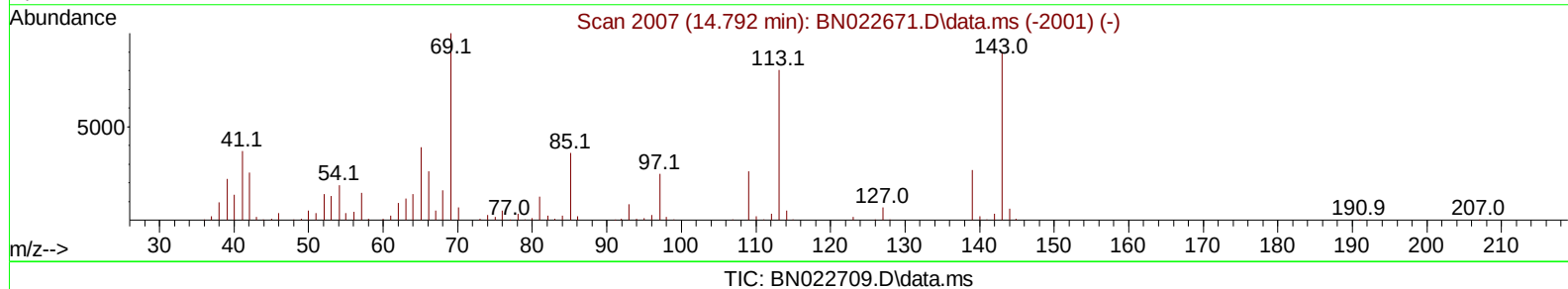
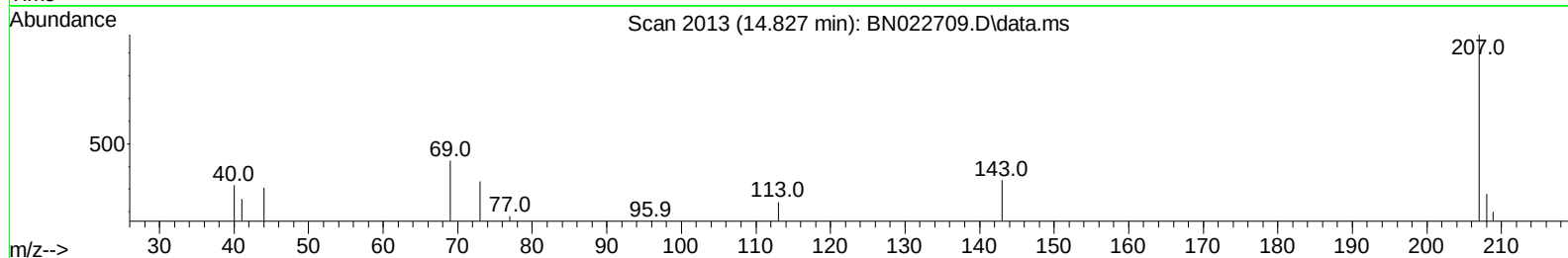
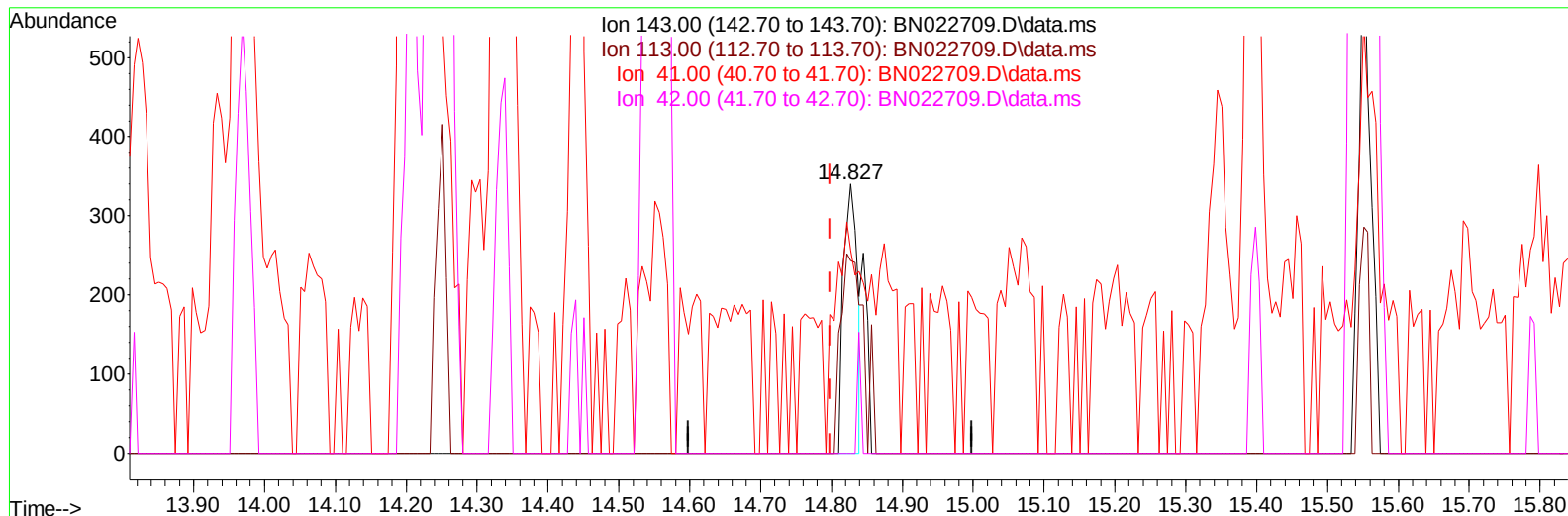
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(54) 4-Nitrophenol-d4 (S)

14.827min (+ 0.029) 0.15 ng/ul

response 472

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	71.47#
41.00	46.30	75.29#
42.00	31.30	0.00#

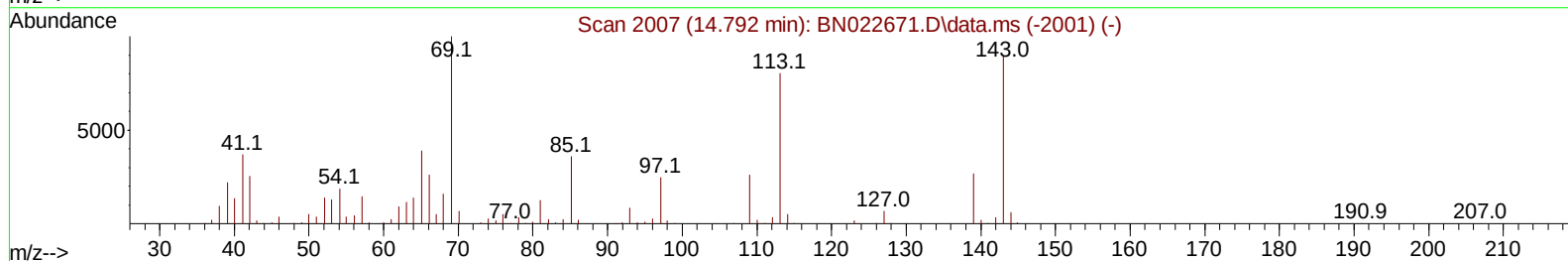
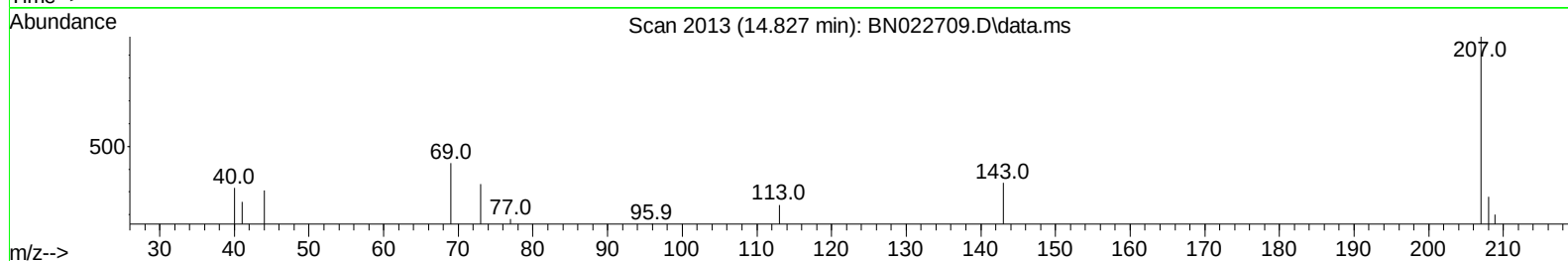
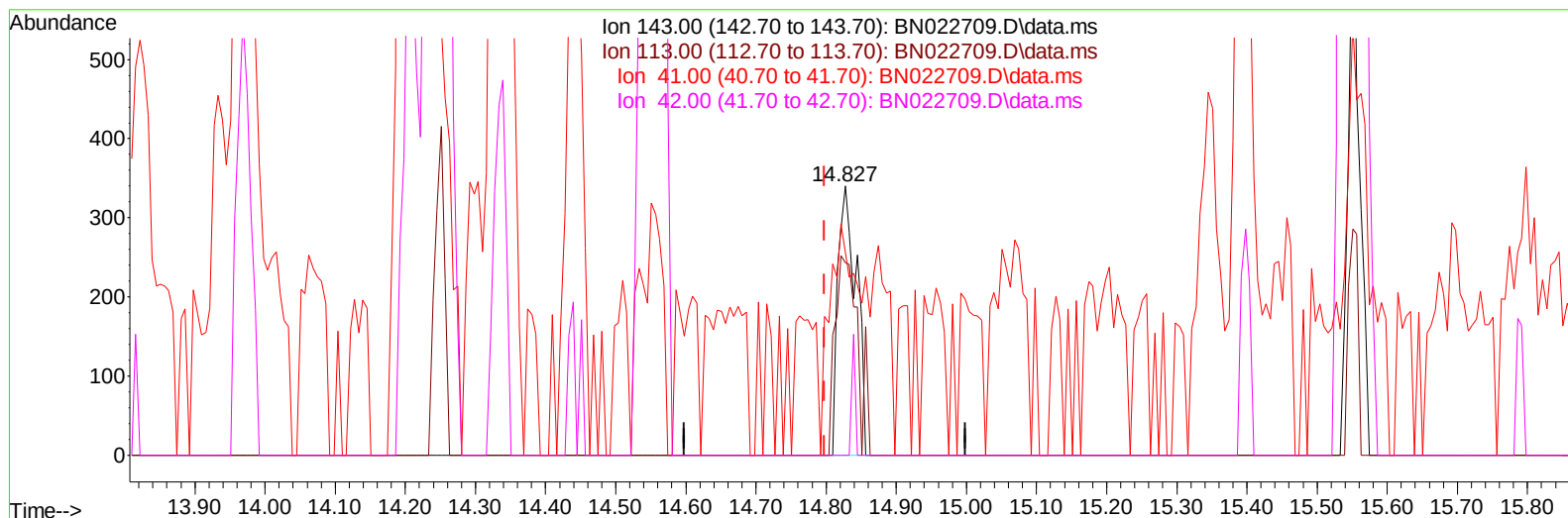
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Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
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TIC: BN022709.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.827min (+ 0.029) 0.19 ng/ul m

response 620

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	71.47#
41.00	46.30	75.29#
42.00	31.30	0.00#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	81086	20.000	ng/ul	0.00
20) Naphthalene-d8	10.692	136	346365	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.557	164	222508	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	441076	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240	397521	20.000	ng/ul	0.00
88) Perylene-d12	23.956	264	367715	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	9969m	4.197	ng/uL	0.00
4) Pyridine-d5	3.634	84	52253	7.780	ng/ul	0.00
7) Phenol-d5	7.040	99	45611	5.956	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.198	67	128949	28.083	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.398	132	107857	19.231	ng/ul	0.00
15) 4-Methylphenol-d8	8.598	113	82579	13.658	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	78650	29.336	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	13163	4.400	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	79799	15.647	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.845	131	147993	19.051	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	493501	30.493	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	515059	28.883	ng/ul	0.00
54) 4-Nitrophenol-d4	14.827	143	620m	0.192	ng/ul	0.03
60) Fluorene-d10	15.551	176	411136	30.974	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	1094	0.390	ng/ul	0.00
73) Anthracene-d10	17.415	188	679342	35.174	ng/ul	0.00
81) Pyrene-d10	19.721	212	732070	35.536	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	645690	34.931	ng/ul	-0.01

Target Compounds Qvalue

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

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