

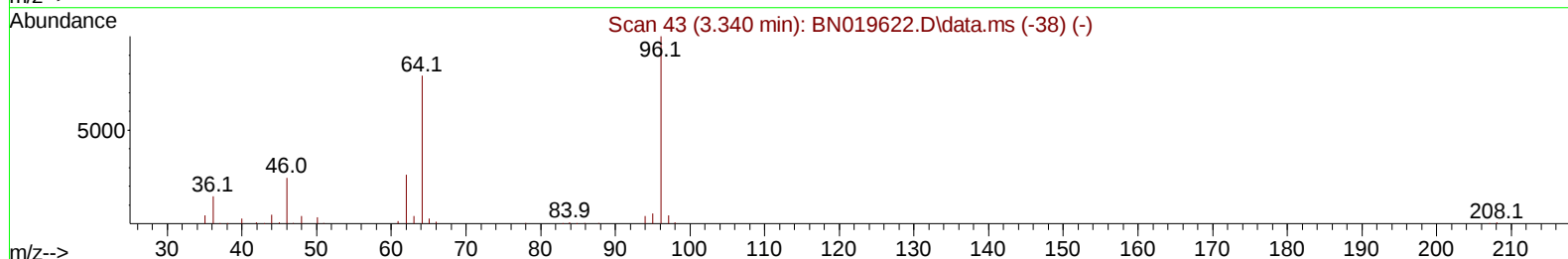
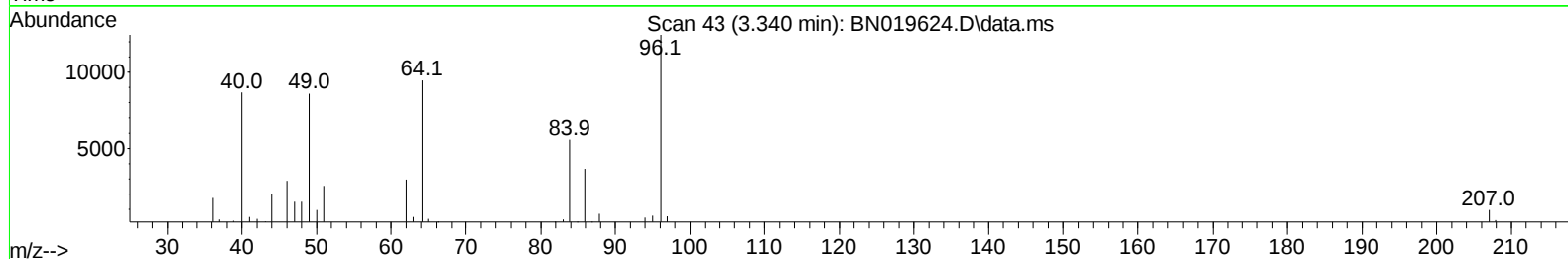
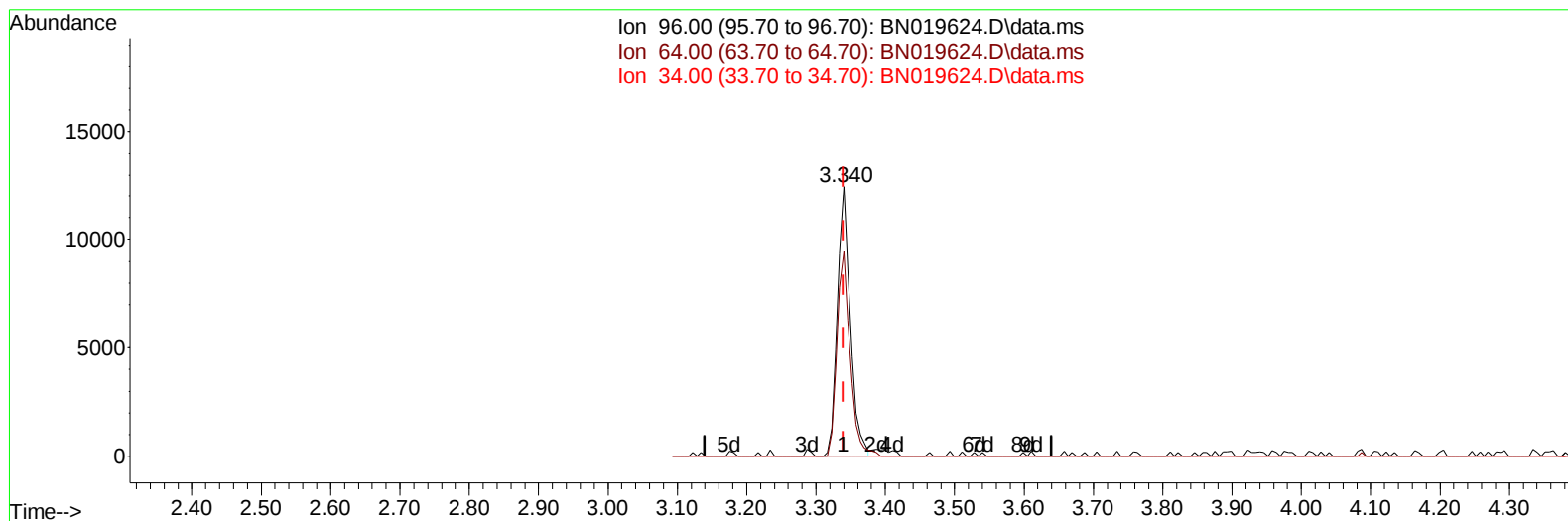
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019624.D
Acq On : 02 May 2022 17:59
Operator : CG/JU
Sample : N2621-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFFJ6

Manual IntegrationsAPPROVED

Quant Time: May 03 03:58:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 03:31:44 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :Yogesh Patel 05/09/2022



TIC: BN019624.D\data.ms

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

3.340min (+ 0.000) 3.99 ng/uL

response 16015

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.08
34.00	0.00	0.00
0.00	0.00	0.00

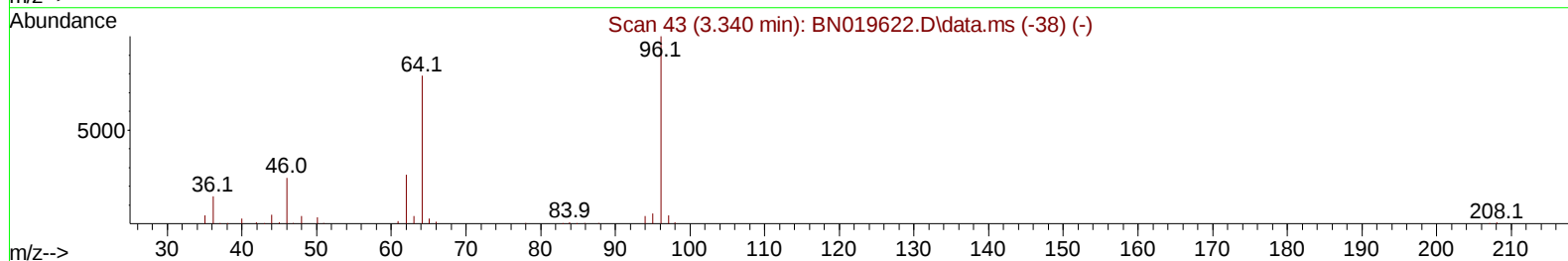
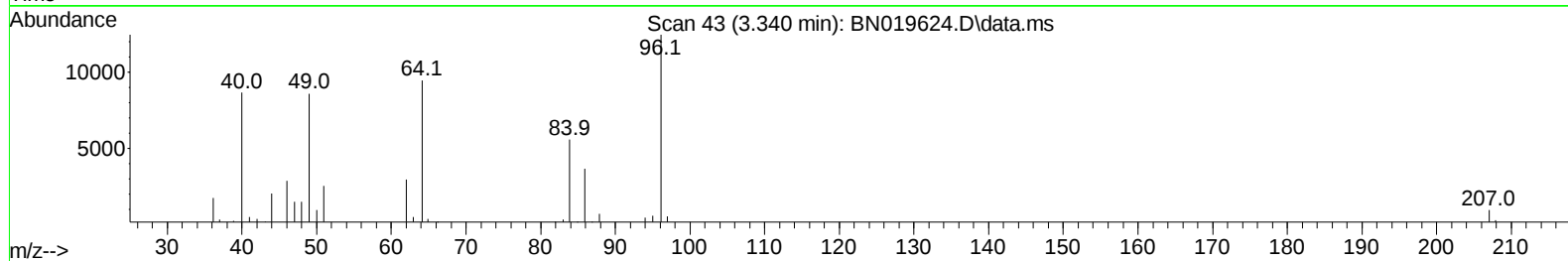
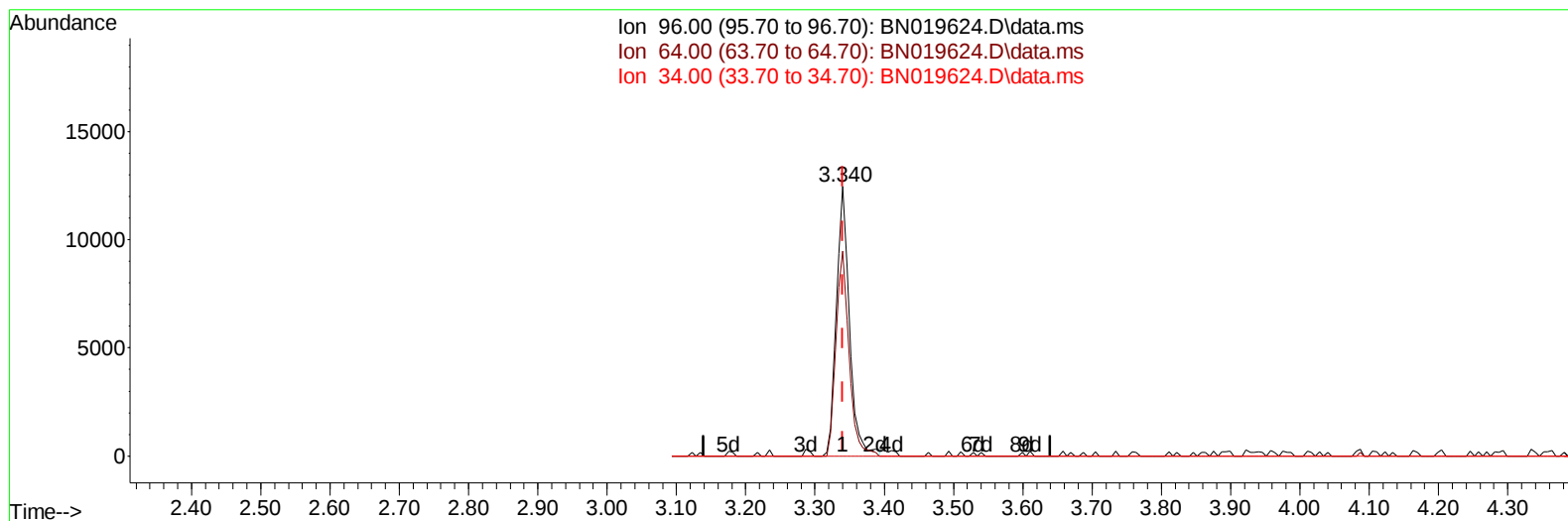
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TIC: BN019624.D\data.ms

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

3.340min (+ 0.000) 4.07 ng/uL m

response 16320

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	76.08
34.00	0.00	0.00
0.00	0.00	0.00

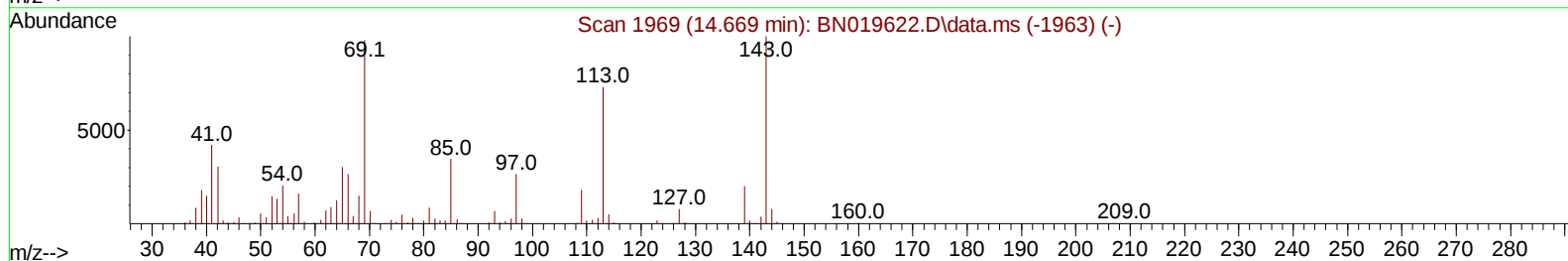
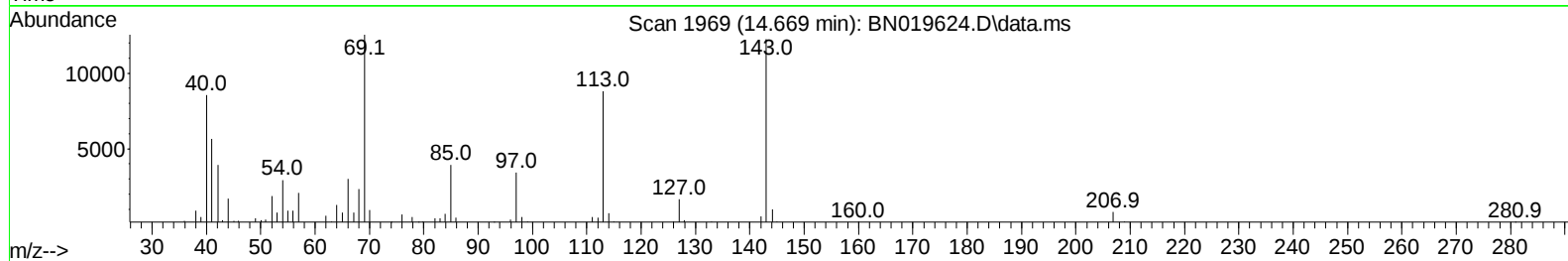
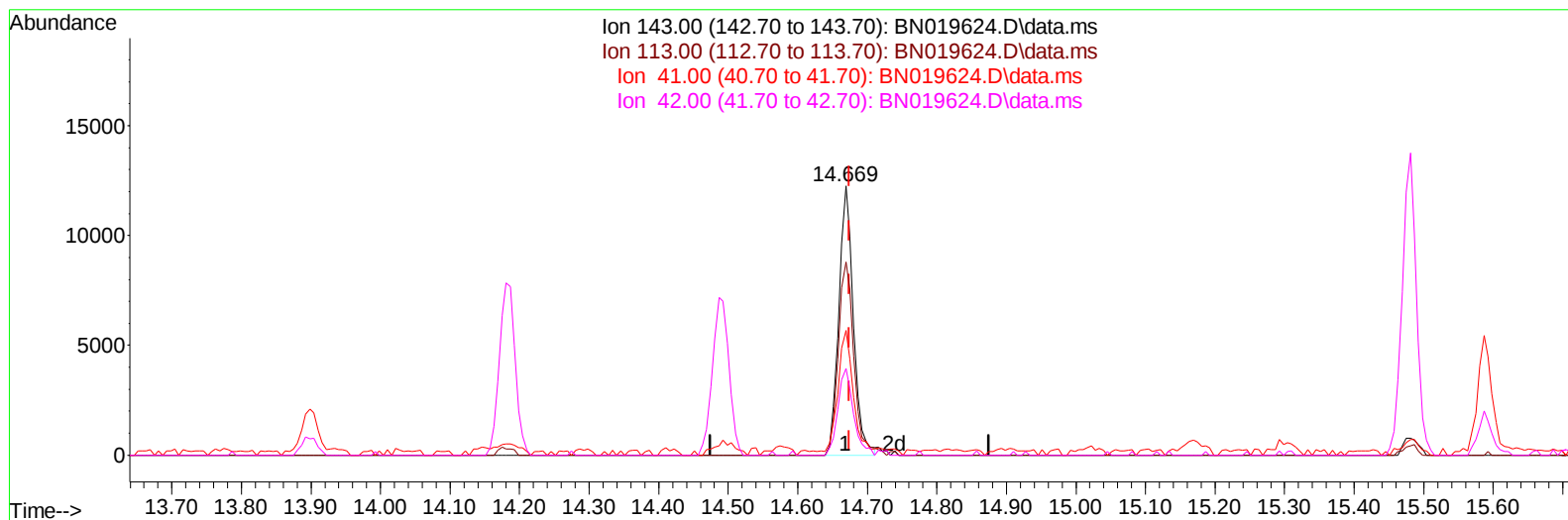
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TIC: BN019624.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 3.74 ng/u1

response 17884

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	71.80
41.00	39.70	46.27
42.00	27.00	32.26

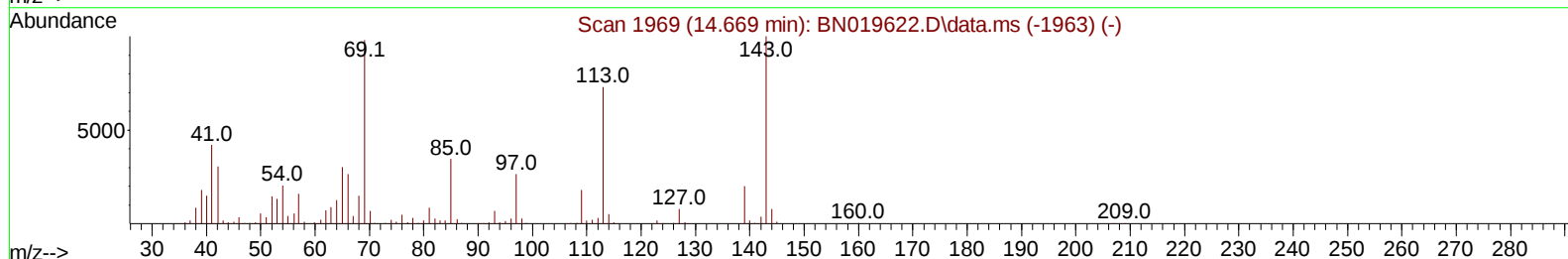
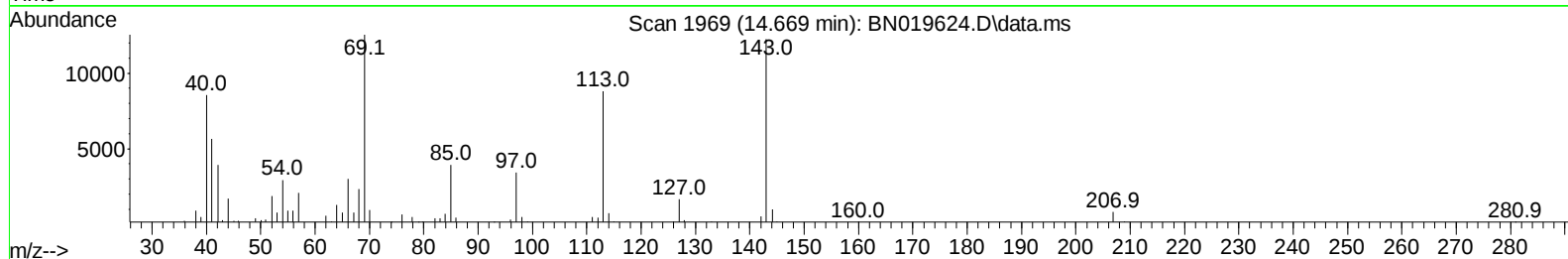
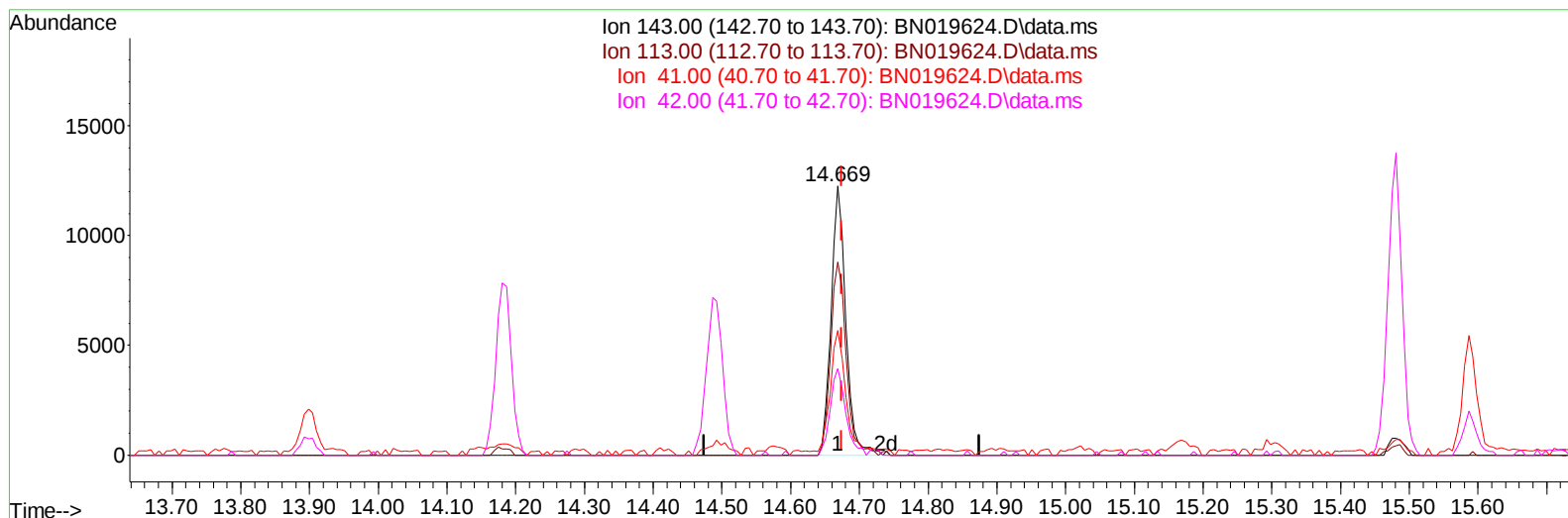
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Operator : CG/JU
Sample : N2621-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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TIC: BN019624.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 3.79 ng/ul m

response 18140

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	71.80
41.00	39.70	46.27
42.00	27.00	32.26

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	151391	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	620256	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	358245	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	691675	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	649769	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	663686	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	16320m	4.068	ng/uL	0.00
4) Pyridine-d5	3.752	84	33848	3.332	ng/ul	0.00
7) Phenol-d5	7.028	99	69916	5.329	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	228281	29.998	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	229158	22.021	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	138617	13.792	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	141187	31.220	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	130038	29.178	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	227704	25.524	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	216189	15.963	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	791514	32.168	ng/ul	0.00
49) Acenaphthylene-d8	14.181	160	963892	28.972	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	18140m	3.791	ng/ul	0.00
60) Fluorene-d10	15.481	176	701143	32.279	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	66123	18.726	ng/ul	0.00
73) Anthracene-d10	17.328	188	1073934	32.970	ng/ul	0.00
81) Pyrene-d10	19.616	212	1173950	31.425	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1124140	33.134	ng/ul	0.00

Target Compounds Qvalue

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

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