

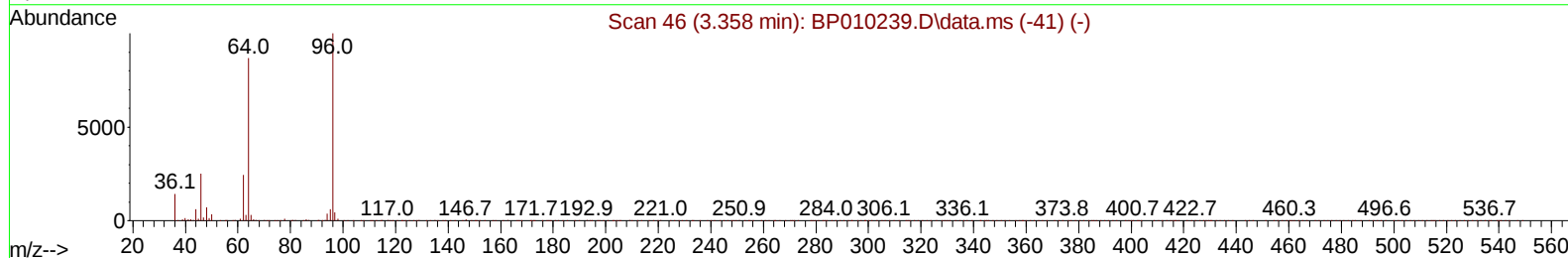
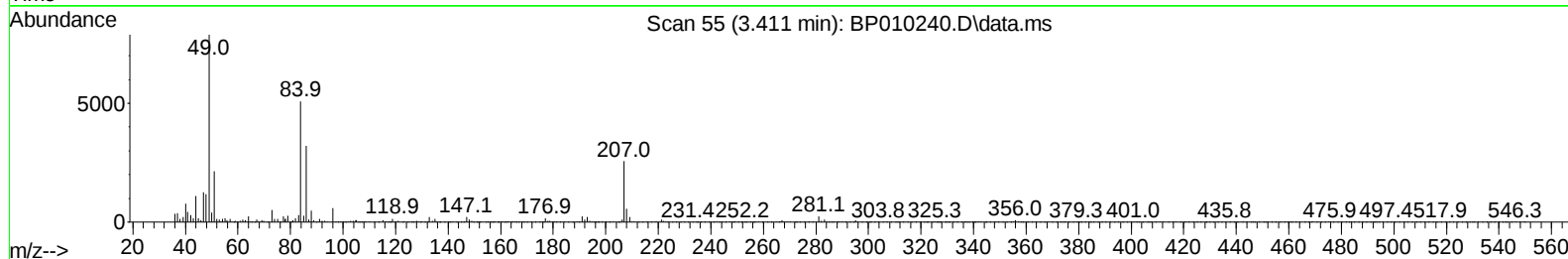
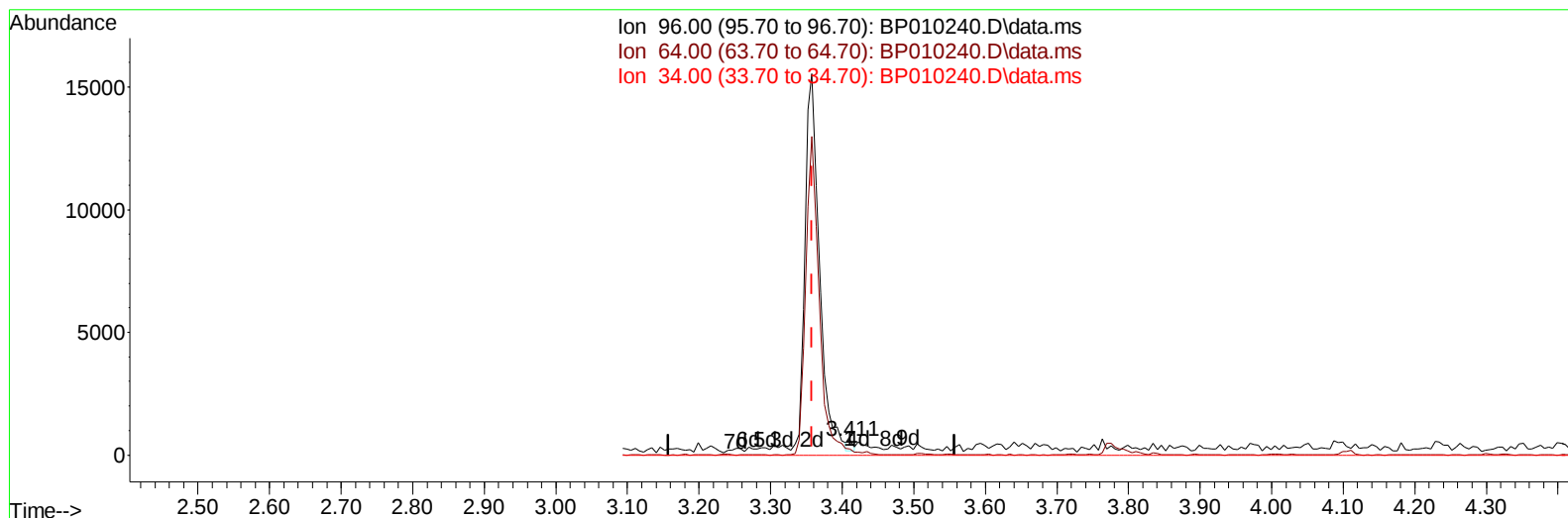
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010240.D
 Acq On : 05 May 2022 11:15
 Operator : CG/JU
 Sample : PB144597BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK597

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/10/2022
 Supervised By :mohammad ahmed 05/10/2022

Quant Time: May 10 08:34:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 06 01:05:07 2022
 Response via : Initial Calibration



TIC: BP010240.D\data.ms

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

3.411min (+ 0.053) 0.05 ng/uL

response 197

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.93
34.00	0.00	0.00
0.00	0.00	0.00

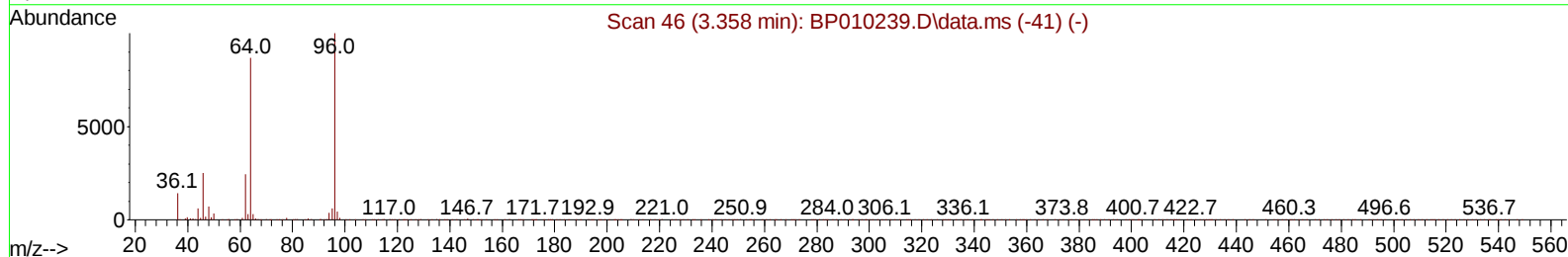
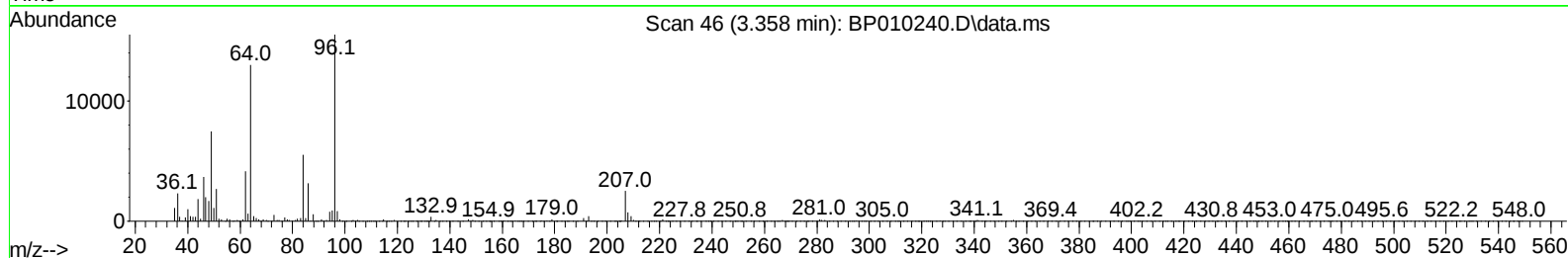
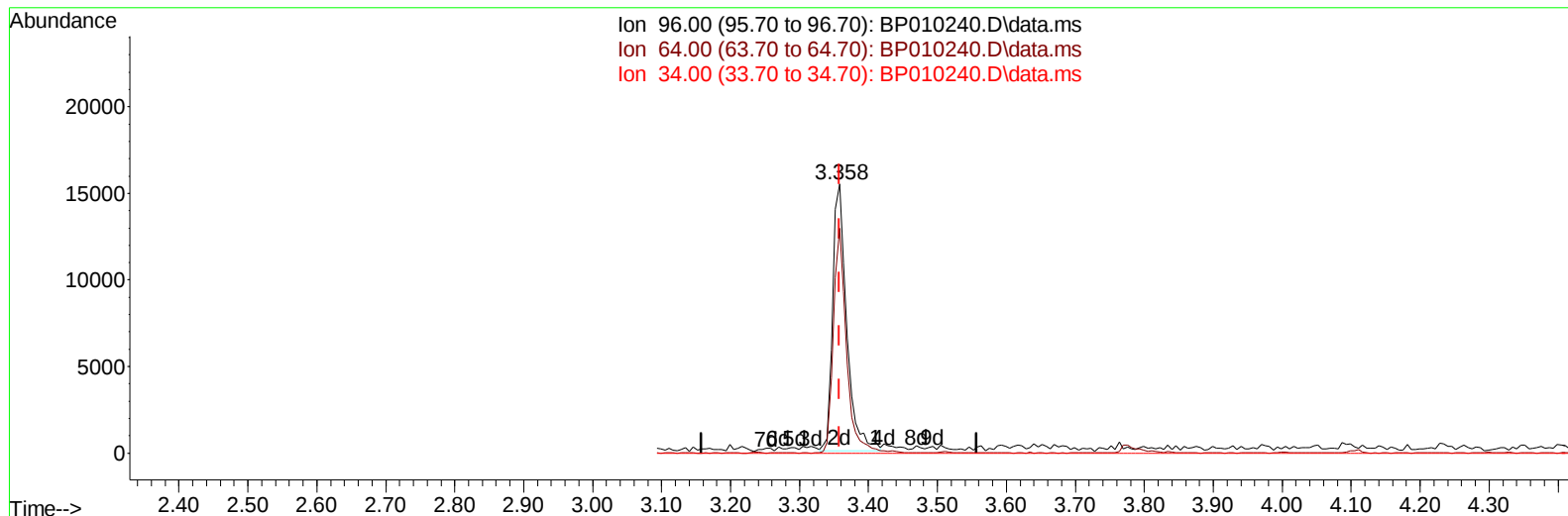
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(3) 1,4-Dioxane-d8 (S)

3.358min (+ 0.000) 5.58 ng/uL m

response 21679

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.58#
34.00	0.00	0.00
0.00	0.00	0.00

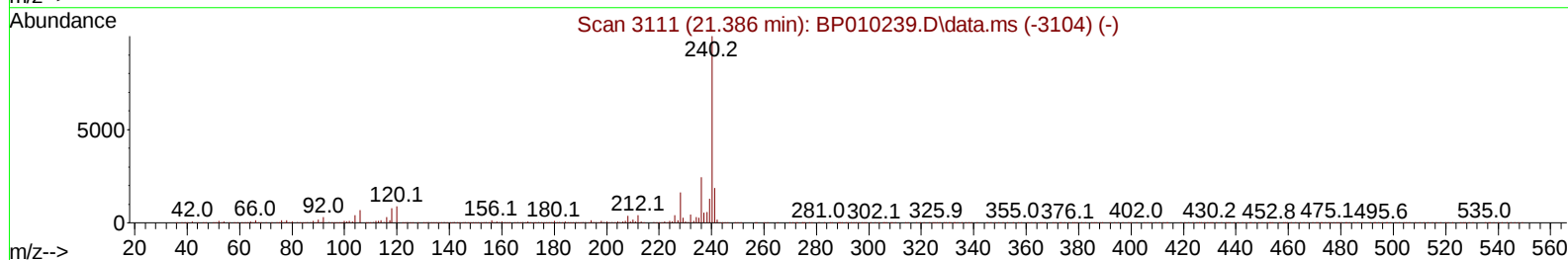
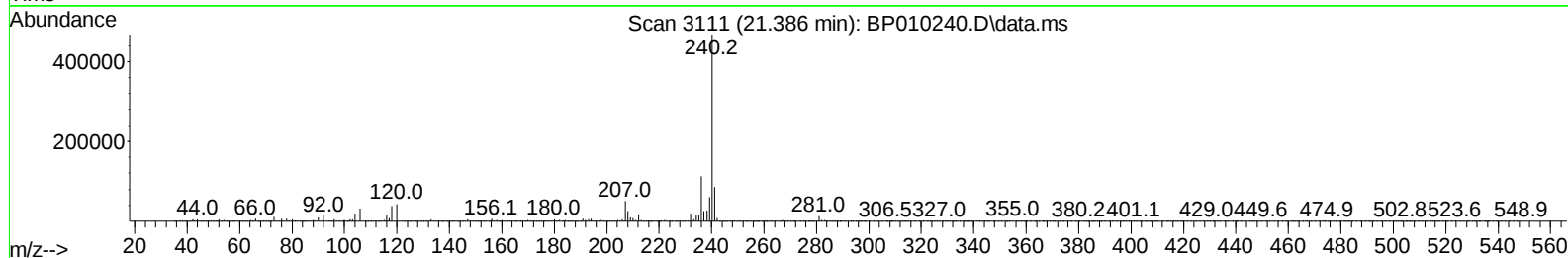
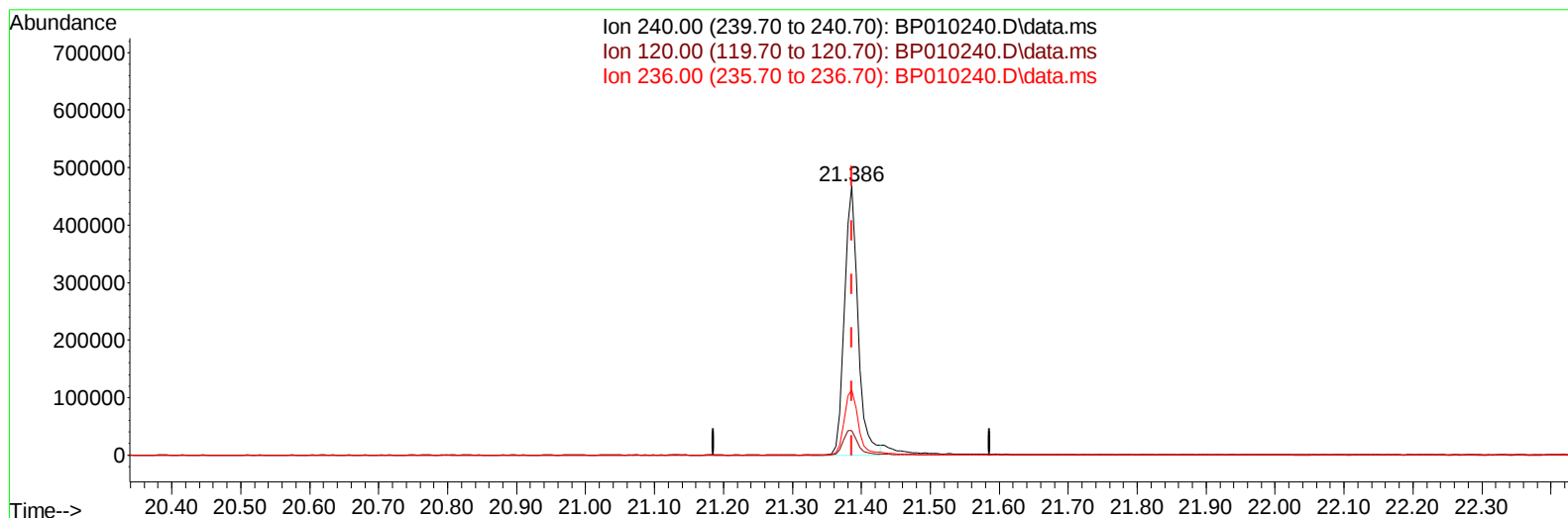
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ALS Vial : 3 Sample Multiplier: 1

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TIC: BP010240.D\data.ms

(79) Chrysene-d12 (I)

21.386min (+ 0.000) 20.00 ng/u1

response 635665

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.22
236.00	31.60	24.11#
0.00	0.00	0.00

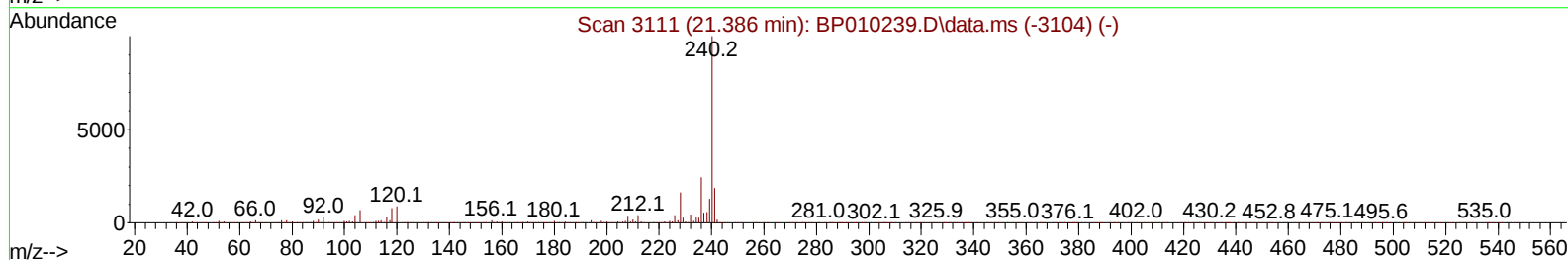
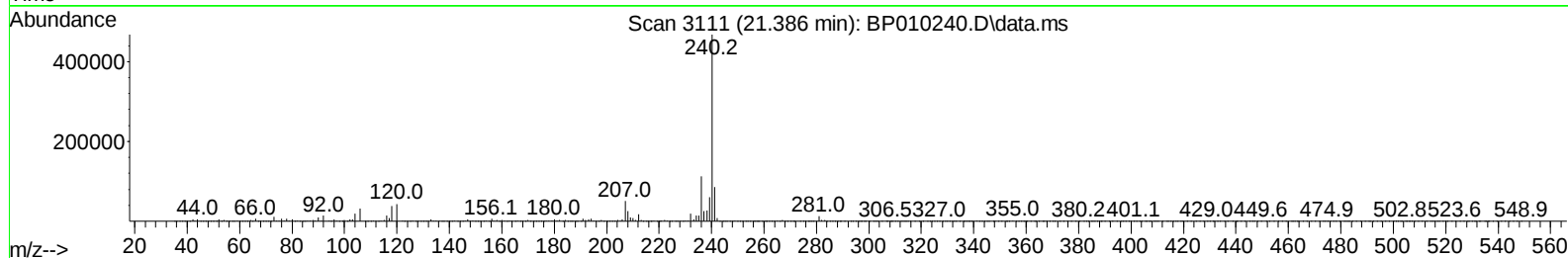
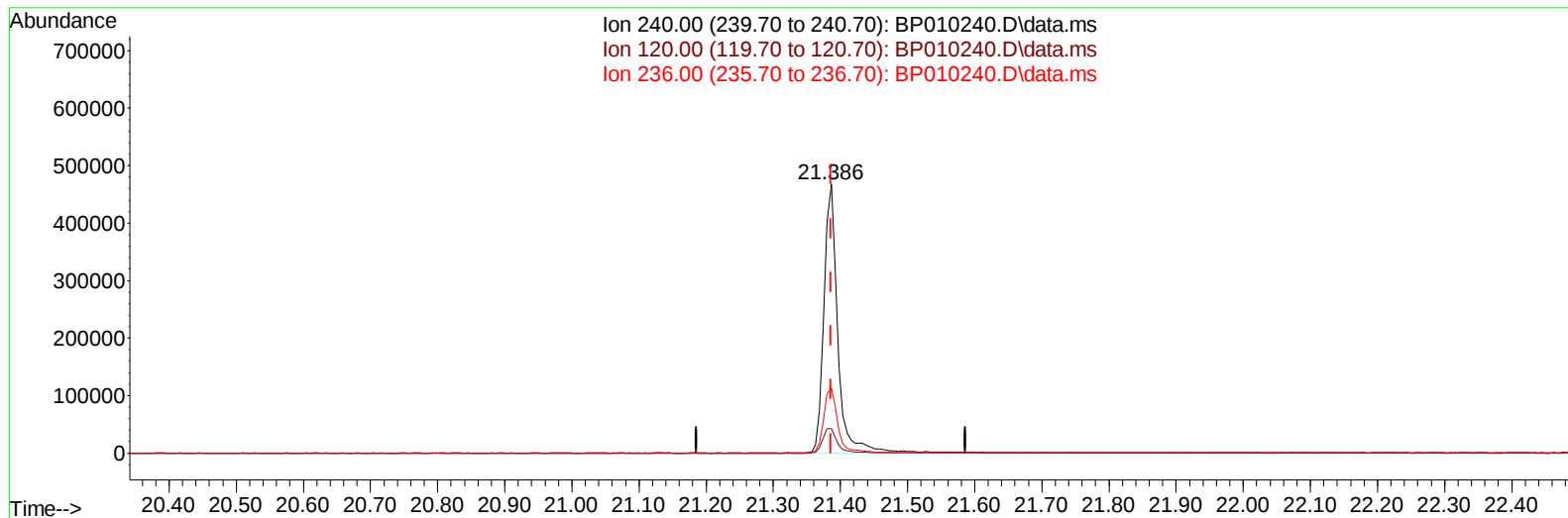
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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TIC: BP010240.D\data.ms

(79) Chrysene-d12 (I)

21.386min (+ 0.000) 20.00 ng/ul m

response 663086

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	9.22
236.00	31.60	24.11#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	169564	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	767622	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	509236	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	950201	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	663086m	20.000	ng/ul	0.00
88) Perylene-d12	23.827	264	613716	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	21679m	5.575	ng/uL	0.00
4) Pyridine-d5	3.775	84	246517	22.535	ng/ul	0.00
7) Phenol-d5	7.081	99	384113	24.904	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	272876	29.485	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	315771	27.645	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	328832	25.733	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	167511	27.437	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	178279	26.737	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	308671	23.780	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	318367	17.106	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1096932	27.691	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1230883	25.674	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	119308	16.360	ng/ul	0.00
60) Fluorene-d10	15.545	176	907537	26.633	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	110215	18.344	ng/ul	0.00
73) Anthracene-d10	17.398	188	1314141	28.587	ng/ul	0.00
81) Pyrene-d10	19.633	212	1291942	34.260	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	965552	30.010	ng/ul	0.00

Target Compounds Qvalue

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

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