

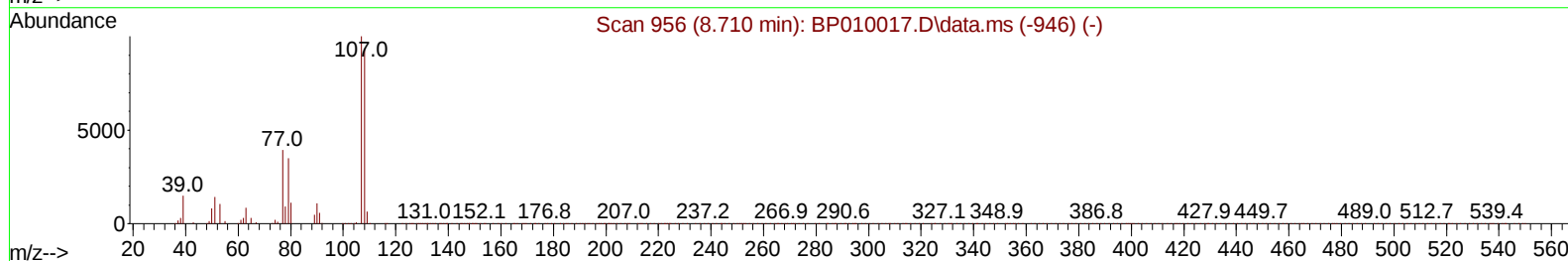
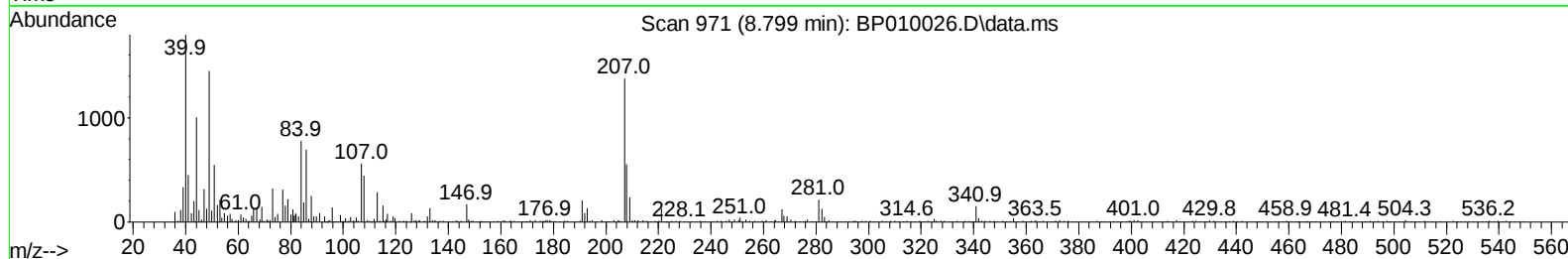
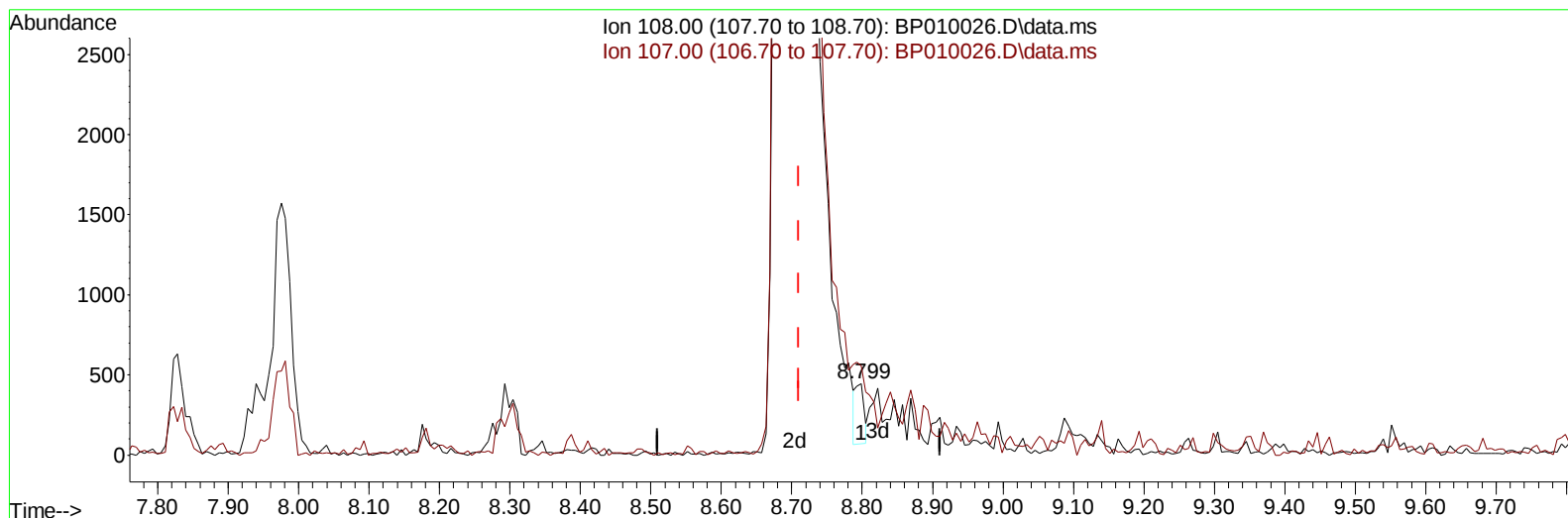
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010026.D
 Acq On : 21 Apr 2022 21:49
 Operator : CG/JU
 Sample : N2473-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J36

Manual IntegrationsAPPROVED

Quant Time: Apr 22 04:20:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010026.D\data.ms

(18) 4-Methylphenol

8.799min (+ 0.088) 0.02 ng/ul

response 306

Ion	Exp%	Act%
108.00	100.00	100.00
107.00	104.50	124.89
0.00	0.00	0.00
0.00	0.00	0.00

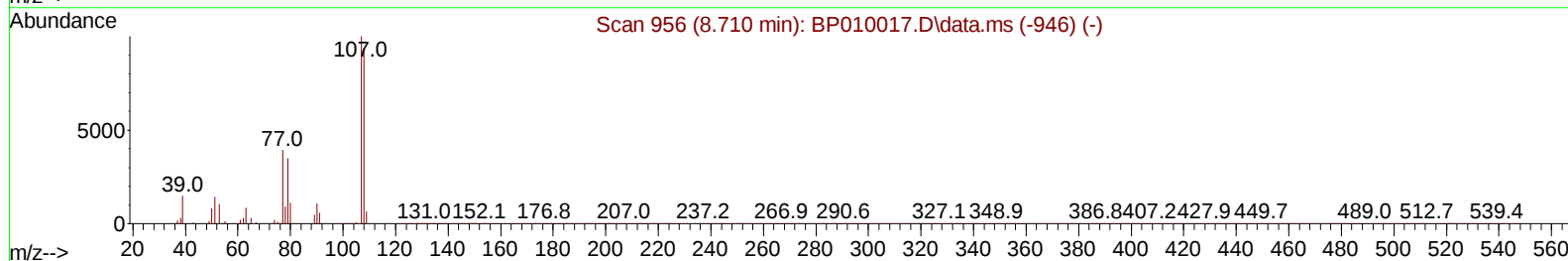
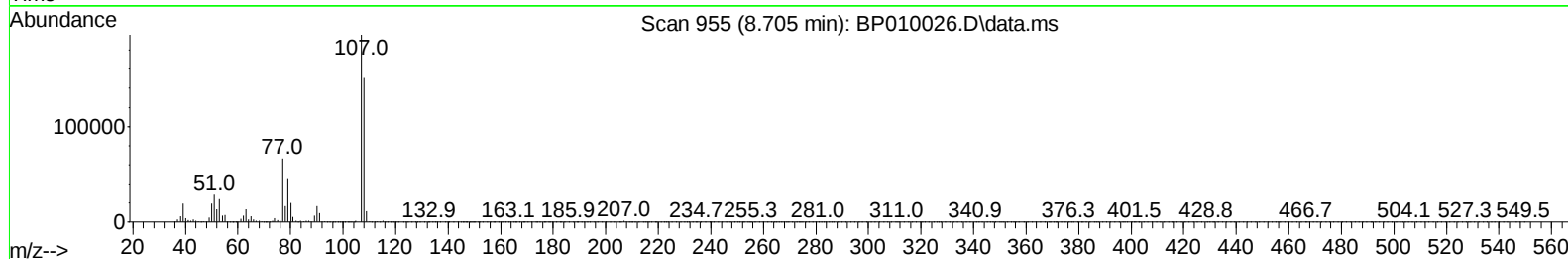
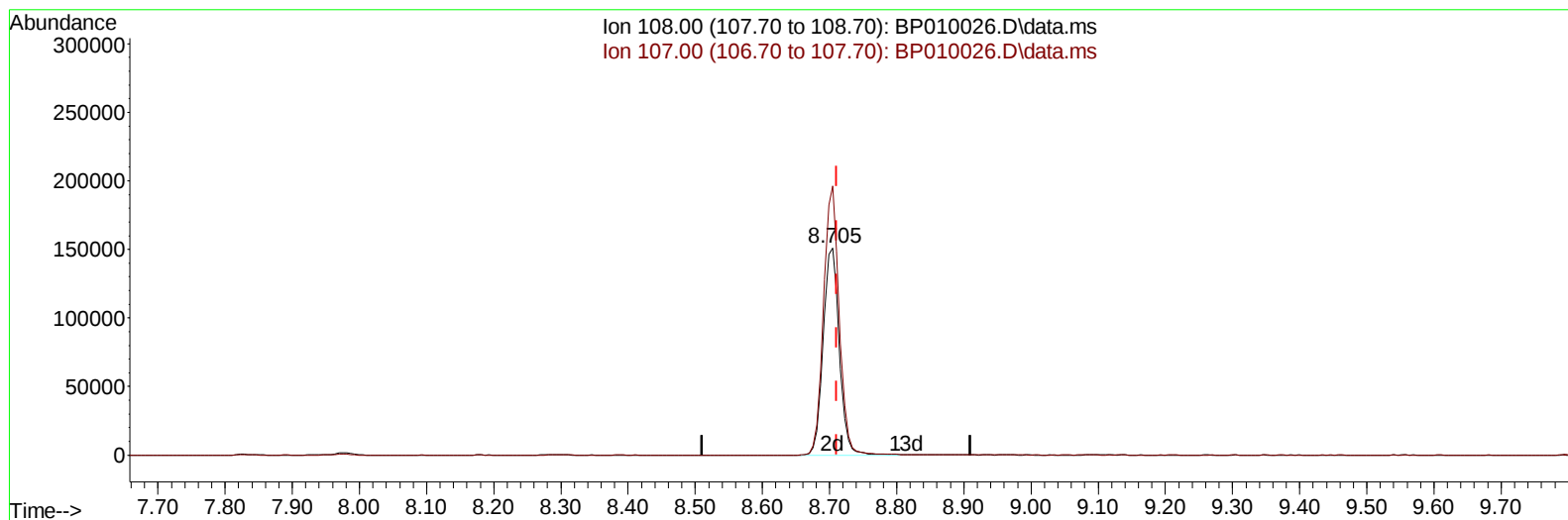
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(18) 4-Methylphenol

8.705min (-0.006) 18.38 ng/ul m

response 251243

Ion	Exp%	Act%
108.00	100.00	100.00
107.00	104.50	129.76#
0.00	0.00	0.00
0.00	0.00	0.00

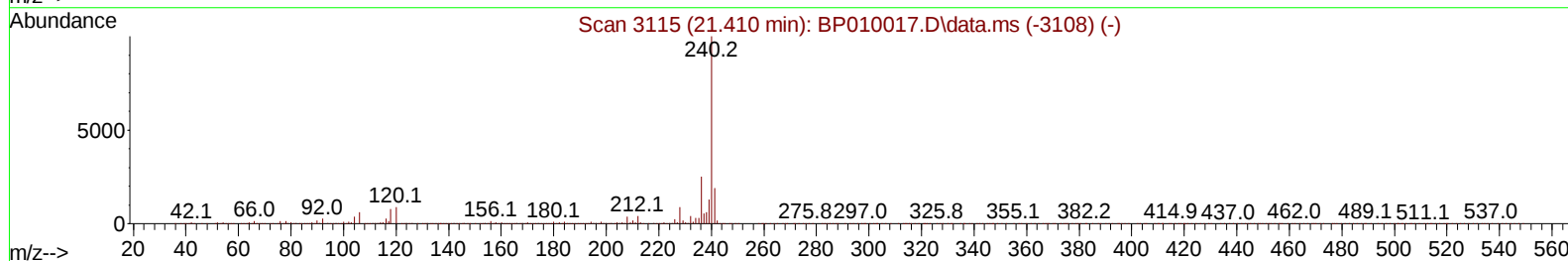
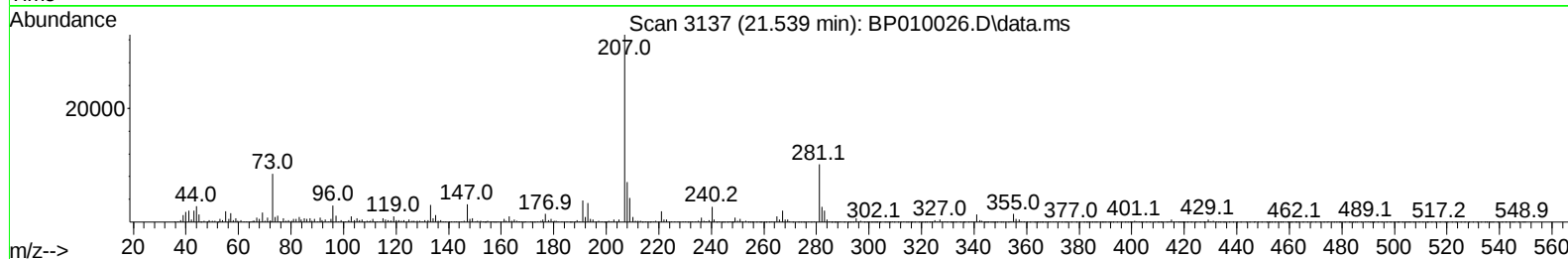
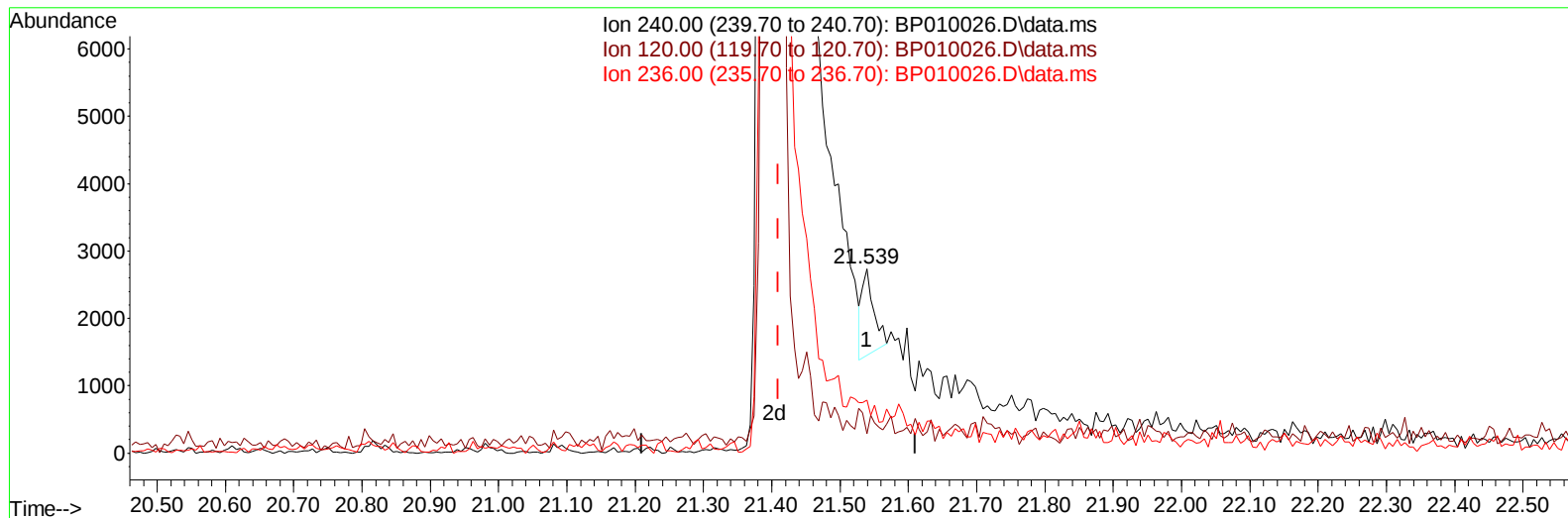
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TIC: BP010026.D\data.ms

(79) Chrysene-d12 (I)

21.539min (+ 0.129) 20.00 ng/u1

response 1544

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	10.67
236.00	31.60	28.61
0.00	0.00	0.00

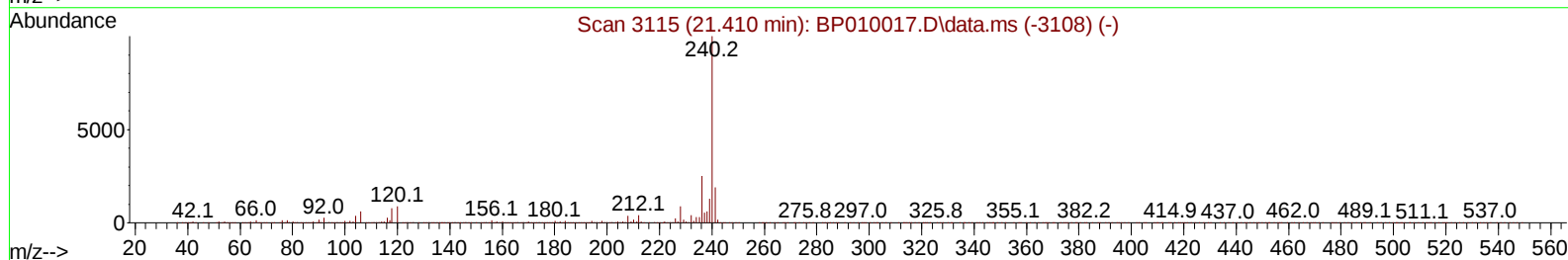
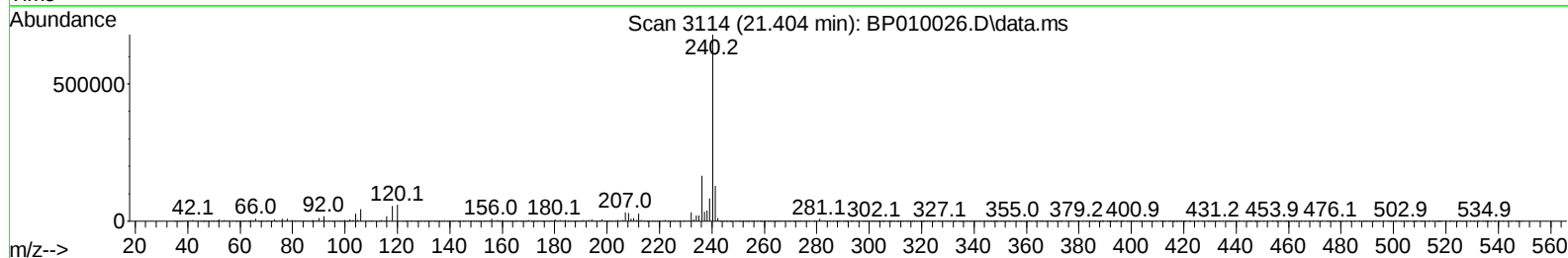
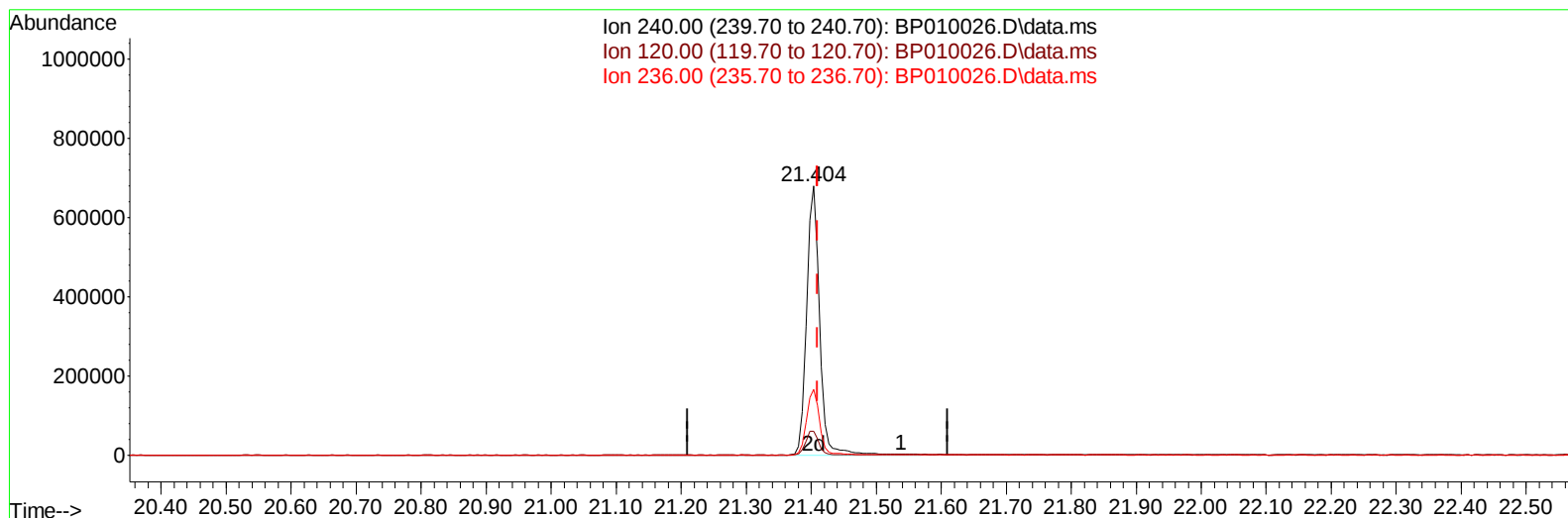
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TIC: BP010026.D\data.ms

(79) Chrysene-d12 (I)

21.404min (-0.006) 20.00 ng/ul m

response 950423

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.78
236.00	31.60	24.51#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	177149	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	806833	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	575336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1217828	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	950423m	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	755250	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	24428	6.013	ng/uL	0.00
4) Pyridine-d5	3.793	84	146553	12.824	ng/ul	0.00
7) Phenol-d5	7.099	99	120187	7.459	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	341441	35.314	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	329315	27.596	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	246283	18.448	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	228216	35.564	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	248091	35.399	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	414117	30.353	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	559711	28.611	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1755346	39.221	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1880296	34.714	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	58863	7.144	ng/ul	0.00
60) Fluorene-d10	15.563	176	1472488	38.248	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	272125	35.339	ng/ul	0.00
73) Anthracene-d10	17.422	188	2363513	40.116	ng/ul	0.00
81) Pyrene-d10	19.651	212	2604678	48.189	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1689116	42.660	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.128	94	111999	6.918	ng/ul#	94
18) 4-Methylphenol	8.705	108	251243m	18.377	ng/ul	

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

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