

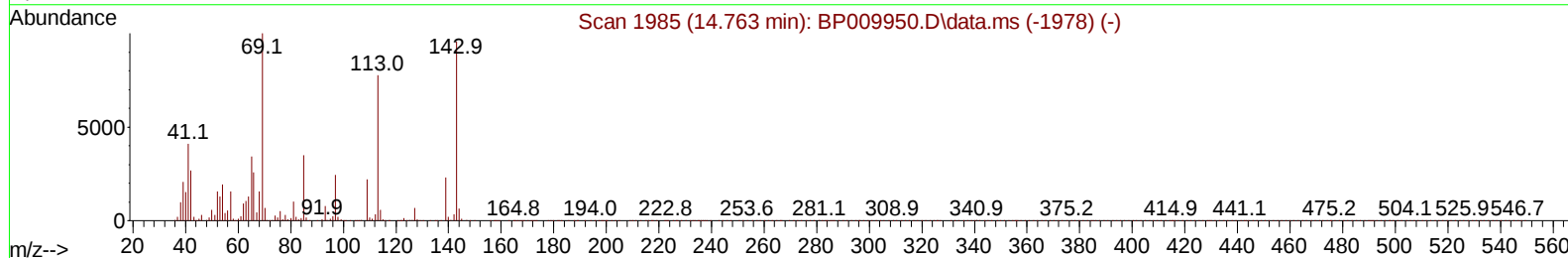
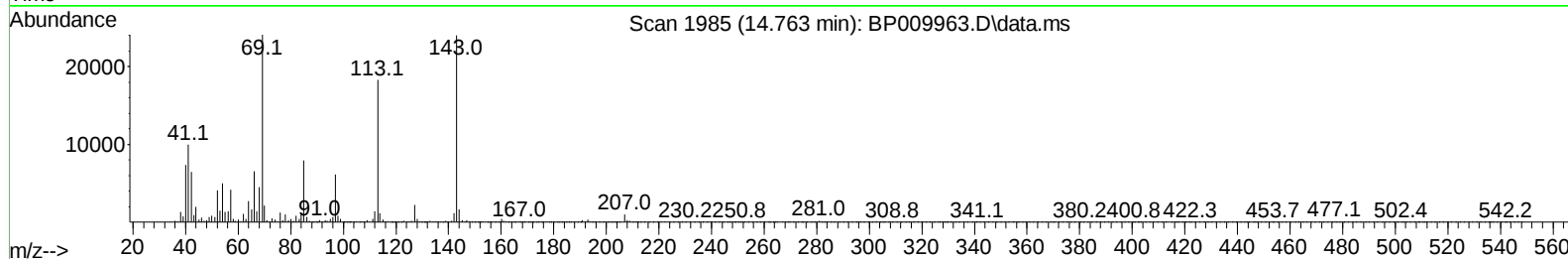
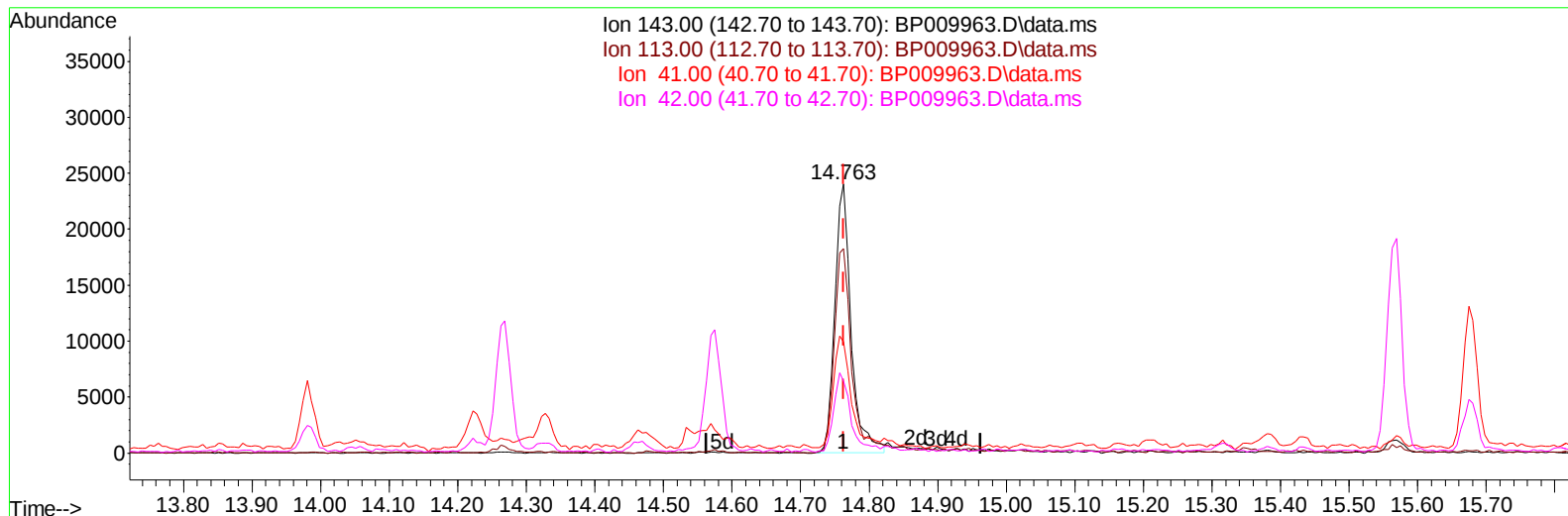
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009963.D
Acq On : 20 Apr 2022 01:21
Operator : CG/JU
Sample : N2422-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J13

Manual IntegrationsAPPROVED

Quant Time: Apr 20 02:38:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 20 01:32:14 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/20/2022
Supervised By :Yogesh Patel 04/25/2022



TIC: BP009963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 5.80 ng/u1

response 39505

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.12#
41.00	23.20	41.46#
42.00	18.00	26.73#

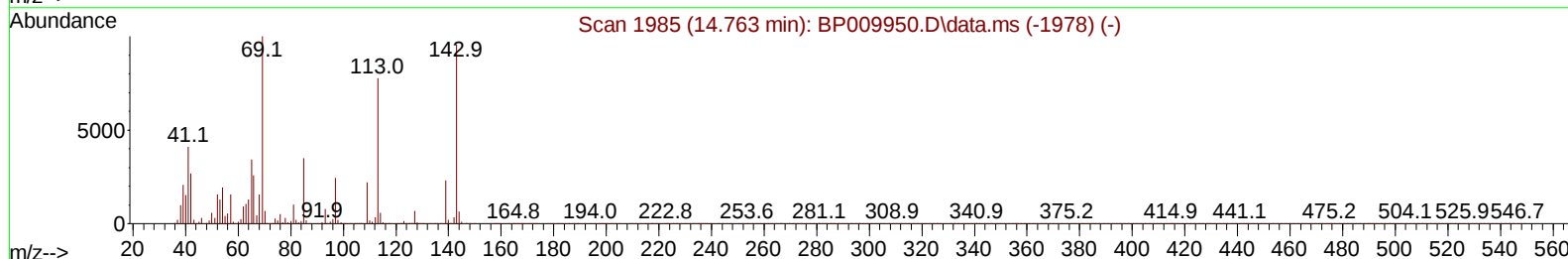
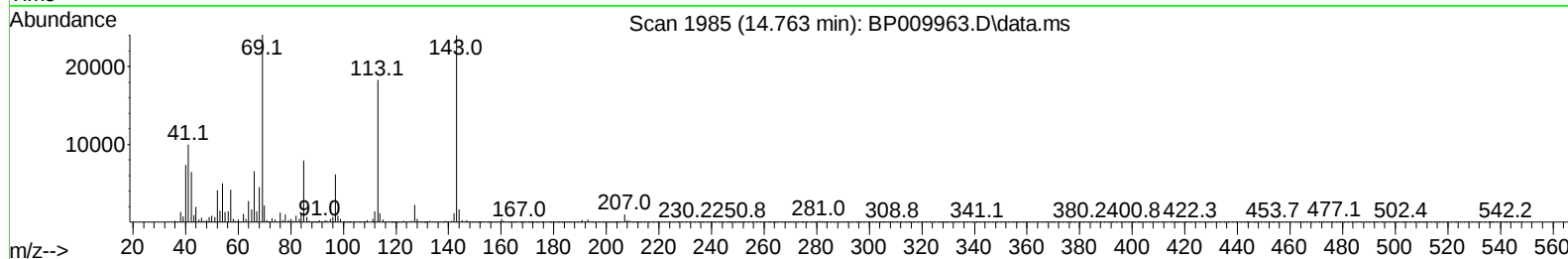
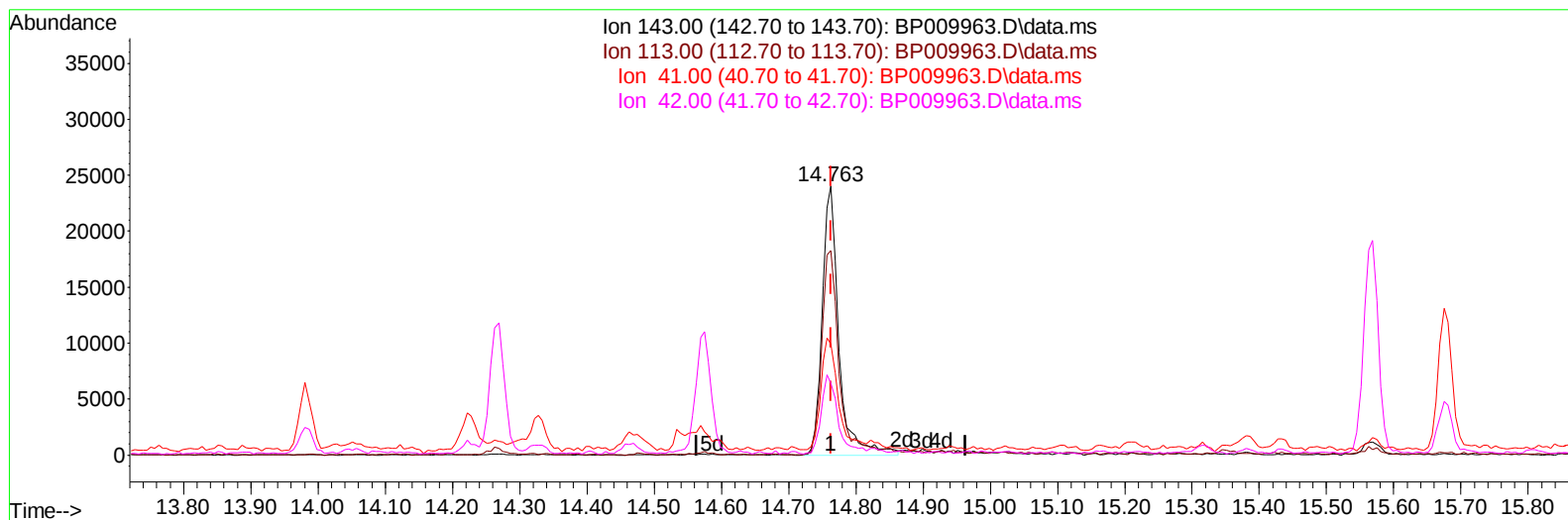
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 Acq On : 20 Apr 2022 01:21
 Operator : CG/JU
 Sample : N2422-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP009963.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 6.01 ng/ul m

response 40939

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.12#
41.00	23.20	41.46#
42.00	18.00	26.73#

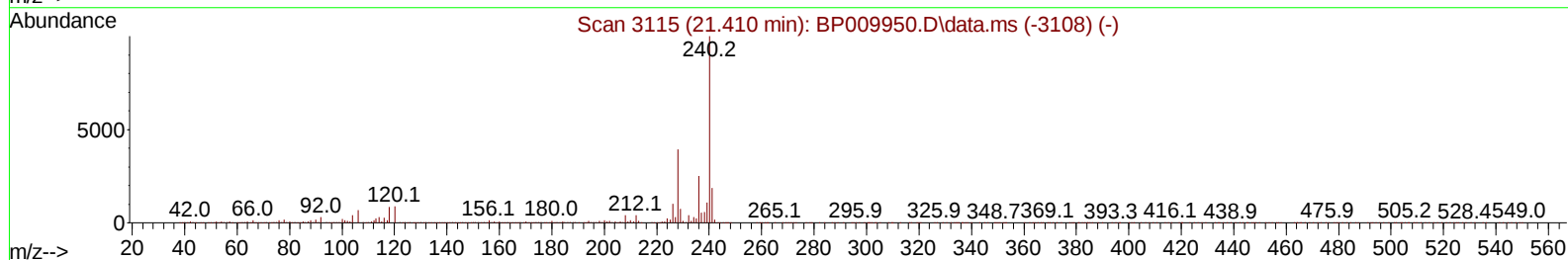
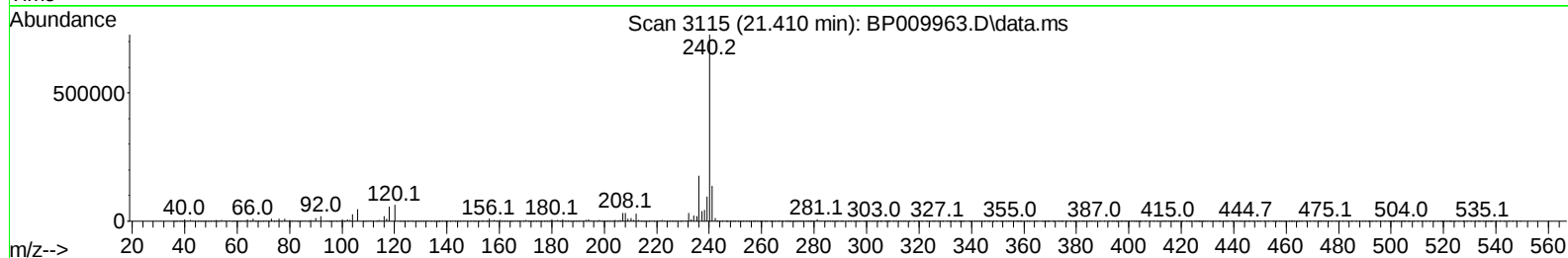
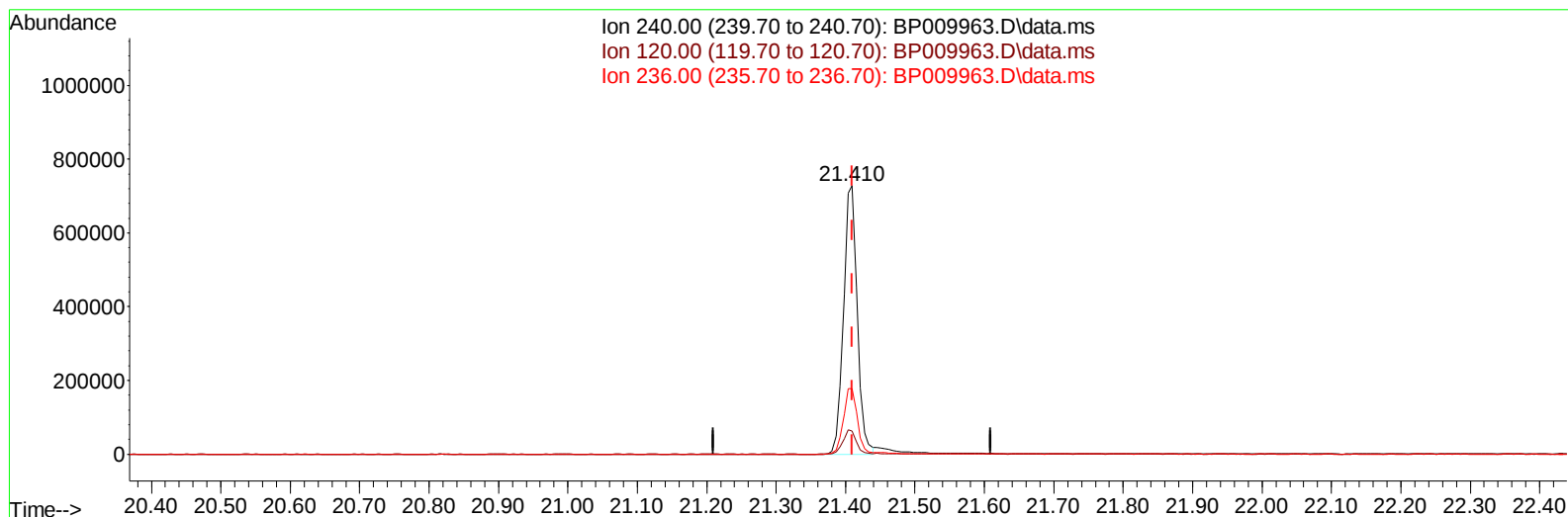
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 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP009963.D\data.ms

(79) Chrysene-d12 (I)

21.410min (-0.000) 20.00 ng/u1

response 1005710

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.71
236.00	31.60	24.34#
0.00	0.00	0.00

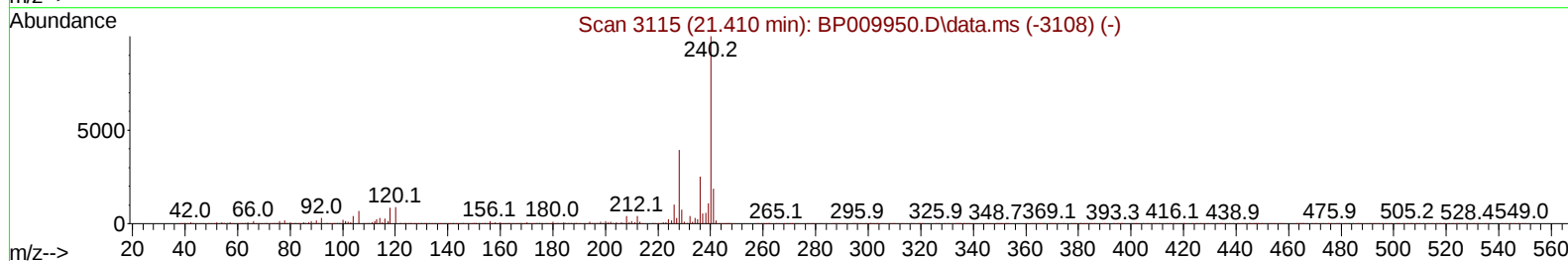
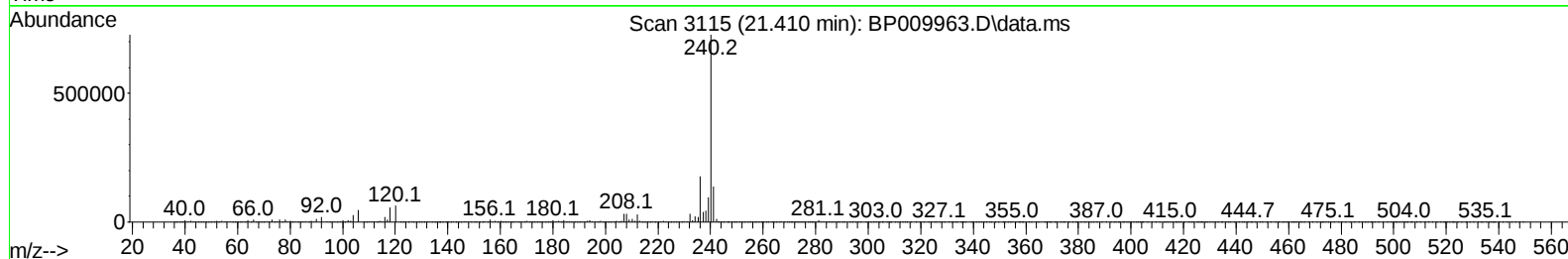
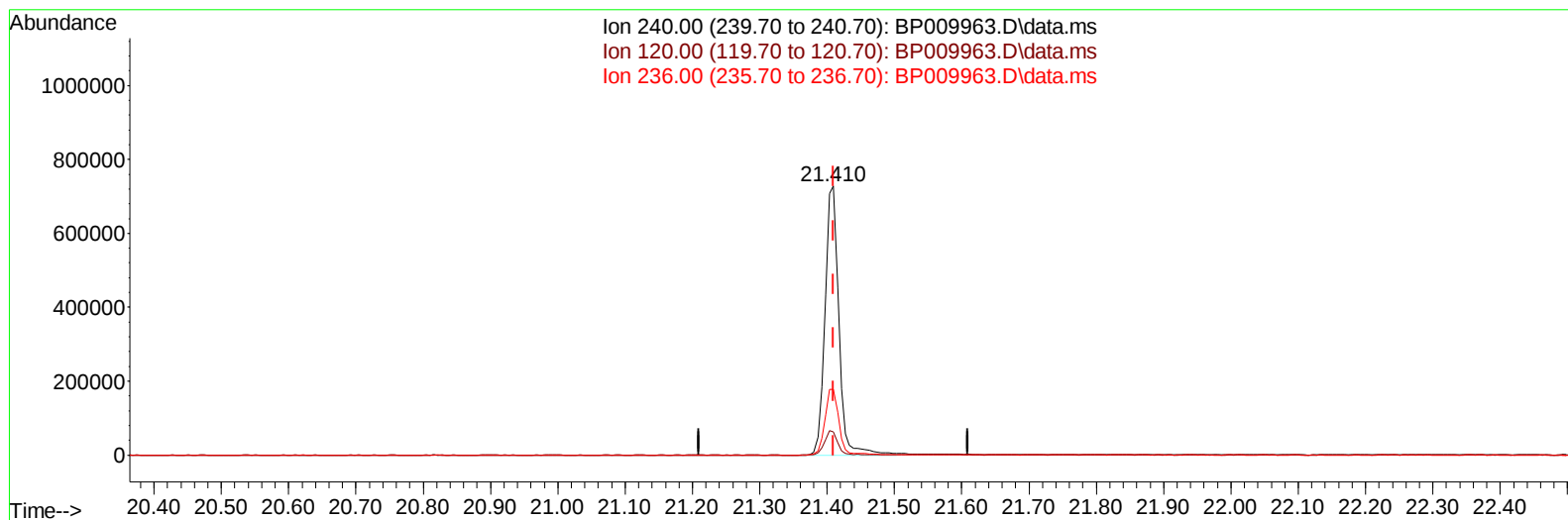
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 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP009963.D\data.ms

(79) Chrysene-d12 (I)

21.410min (-0.000) 20.00 ng/ul m

response 1045046

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.71
236.00	31.60	24.34#
0.00	0.00	0.00

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 ALS Vial : 15 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	139556	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	643381	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	478131	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1054971	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	1045046m	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	963093	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	16137	5.161	ng/uL	0.00
4) Pyridine-d5	3.793	84	64679	7.502	ng/ul	0.00
7) Phenol-d5	7.105	99	80755	6.591	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	235785	32.325	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	223725	24.208	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	165373	16.169	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	151952	32.751	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	164028	32.247	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	285317	26.805	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	381217	25.146	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1263441	33.640	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1346598	30.264	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	40939m	6.010	ng/ul	0.00
60) Fluorene-d10	15.569	176	1077253	34.096	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	209798	33.569	ng/ul	0.00
73) Anthracene-d10	17.428	188	1874369	37.108	ng/ul	0.00
81) Pyrene-d10	19.657	212	2296614	40.238	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1942380	38.728	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.128	94	14246	1.162	ng/ul#	85

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

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