Quantitation Report (Qedit)

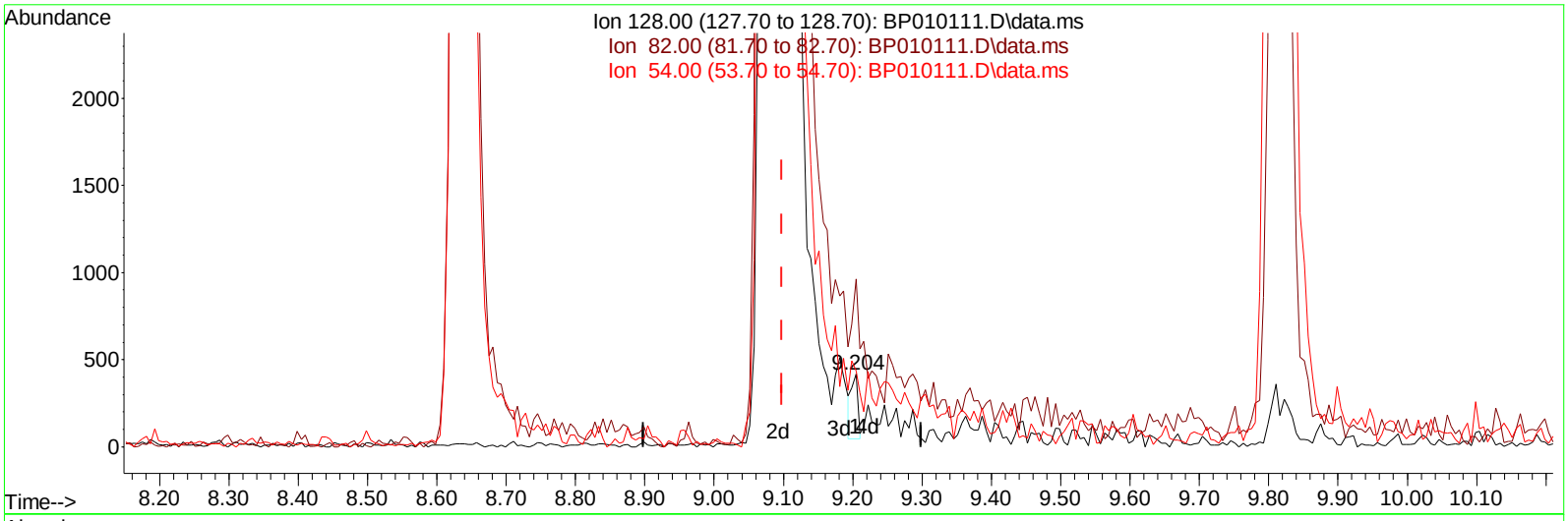
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010111.D
Acq On : 26 Apr 2022 12:25
Operator : CG/JU
Sample : N2517-01
Misc :
ALS Vial : 33 Sample Multiplier: 1

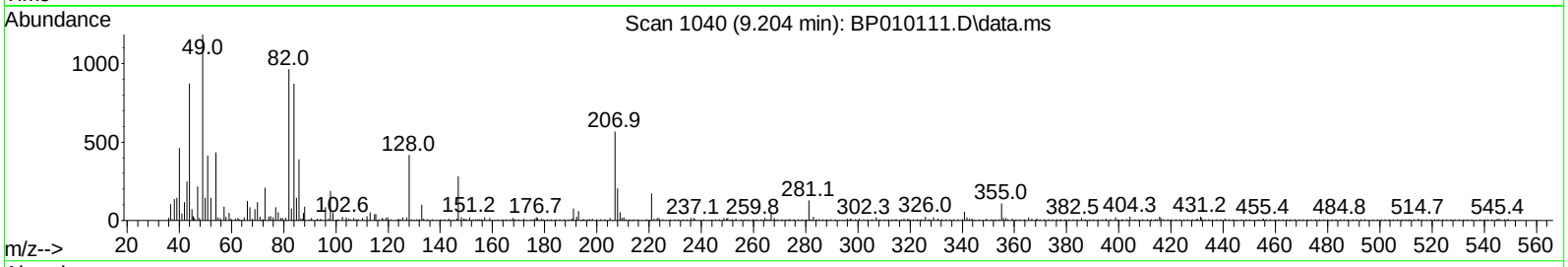
Quant Time: Apr 27 00:44:58 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

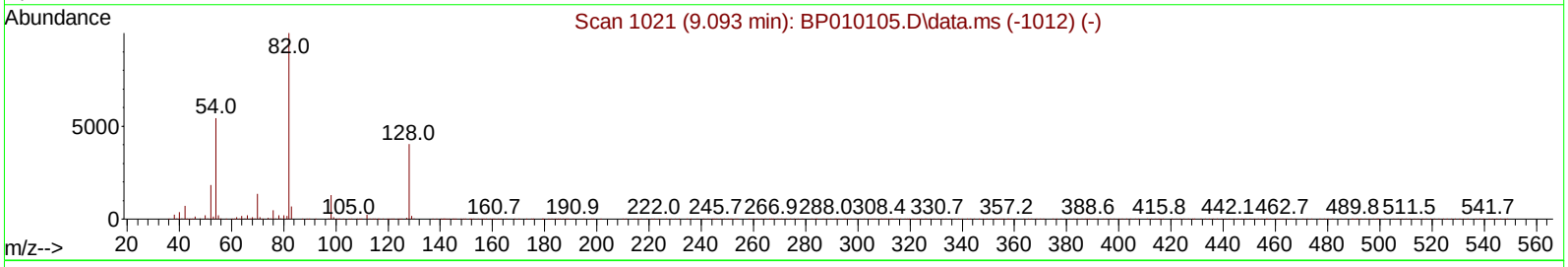
Instrument :
BNA_P
ClientSampleId :
BFFH1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/27/2022
Supervised By :Yogesh Patel 04/28/2022







TIC: BP010111.D\data.ms

(21) Nitrobenzene-d5 (S)

9.204min (+ 0.106) 0.06 ng/ul

response 245

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	231.18
54.00	120.20	109.35
0.00	0.00	0.00

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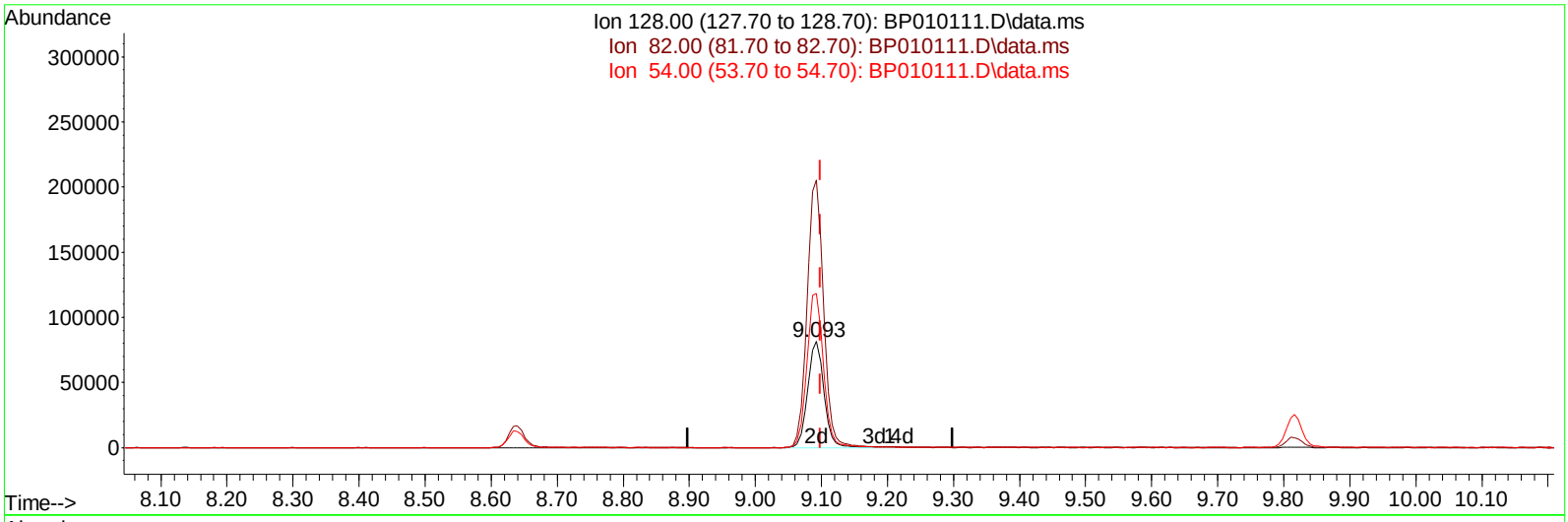

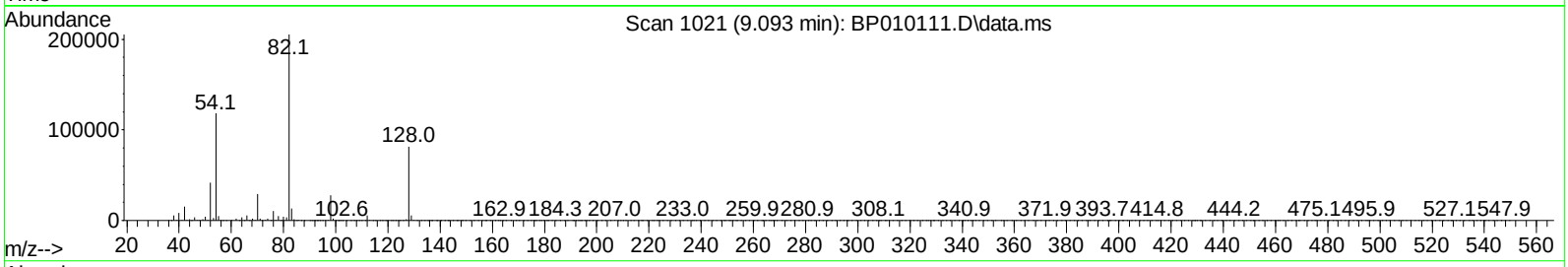

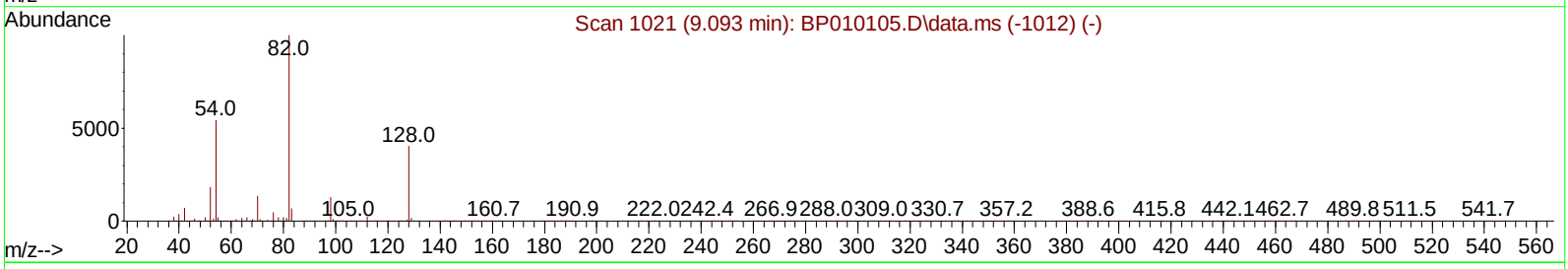

TIC: BP010111.D\data.ms

(21) Nitrobenzene-d5 (S)

9.093min (-0.006) 31.52 ng/ul m

response 138913

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	252.13
54.00	120.20	145.16#
0.00	0.00	0.00

Quantitation Report (Qedit)

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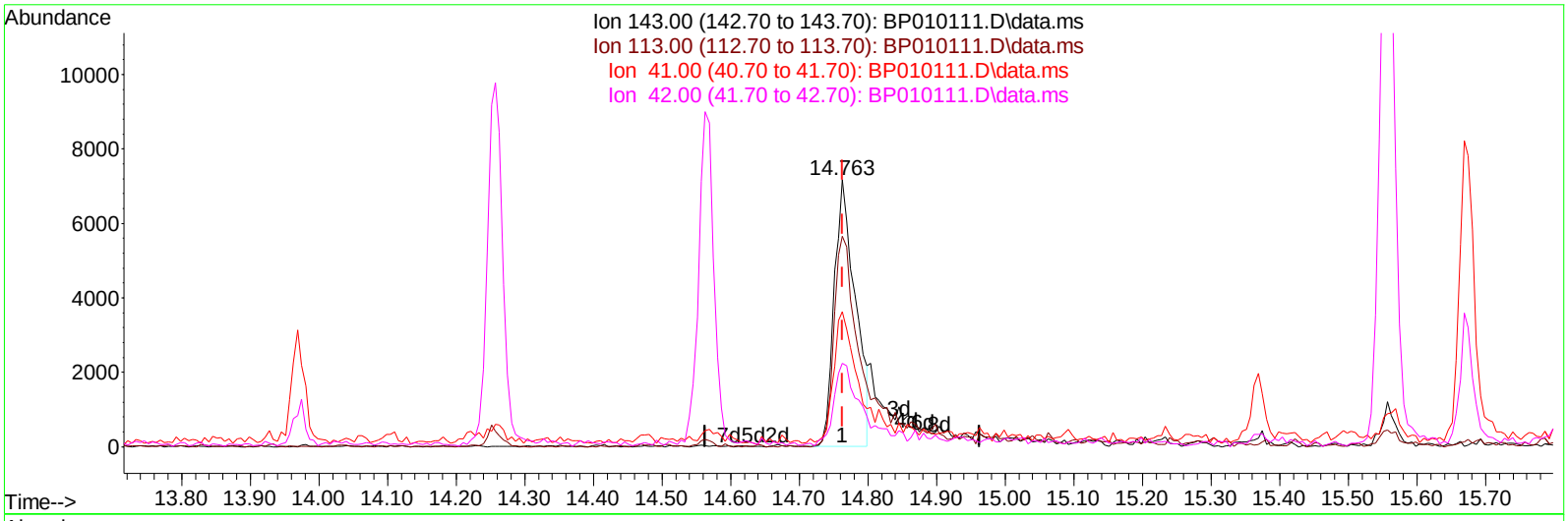

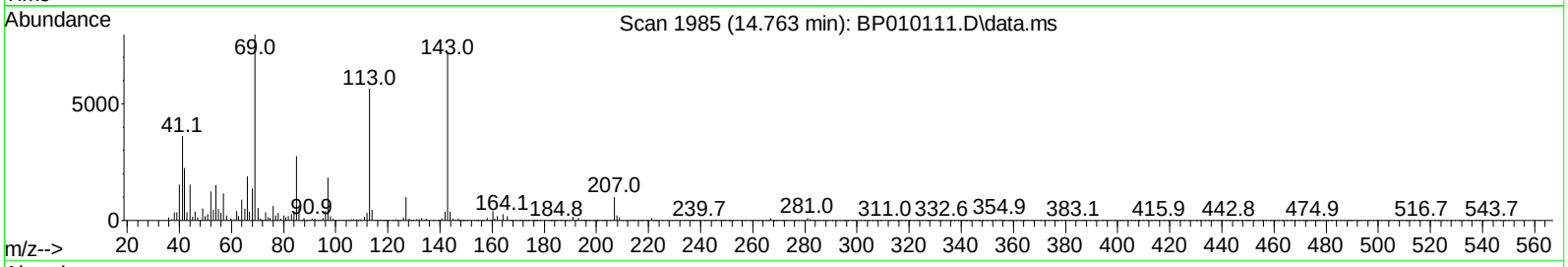

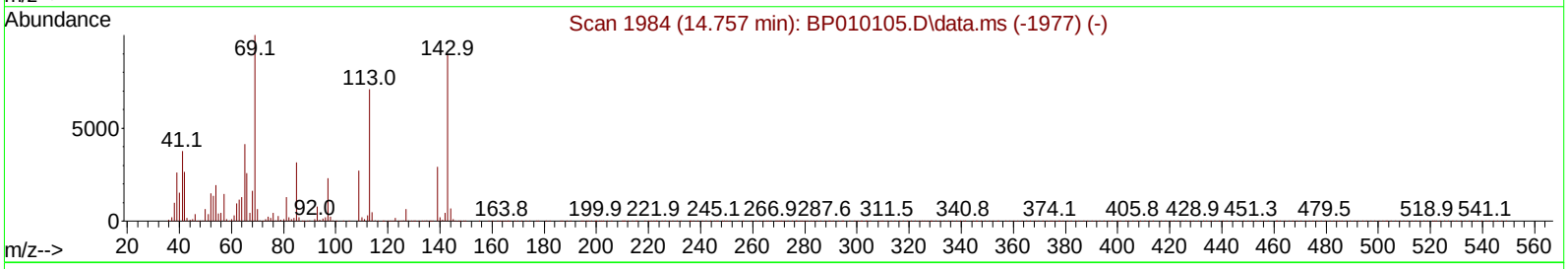

TIC: BP010111.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 2.89 ng/ul

response 15379

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.77#
41.00	23.20	50.57#
42.00	18.00	31.37#

Quantitation Report (Qedit)

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Operator : CG/JU
Sample : N2517-01
Misc :
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BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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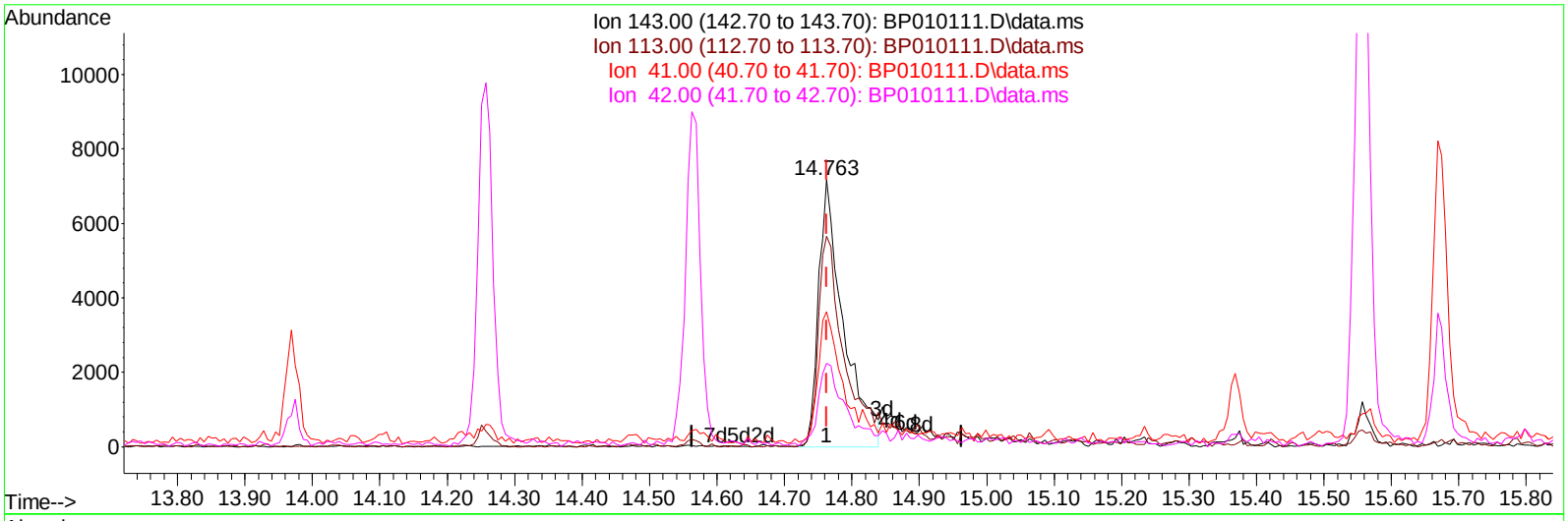

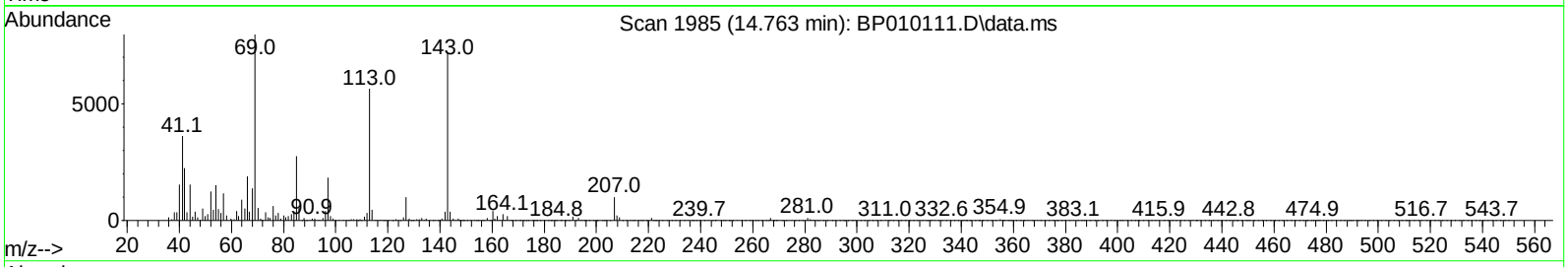

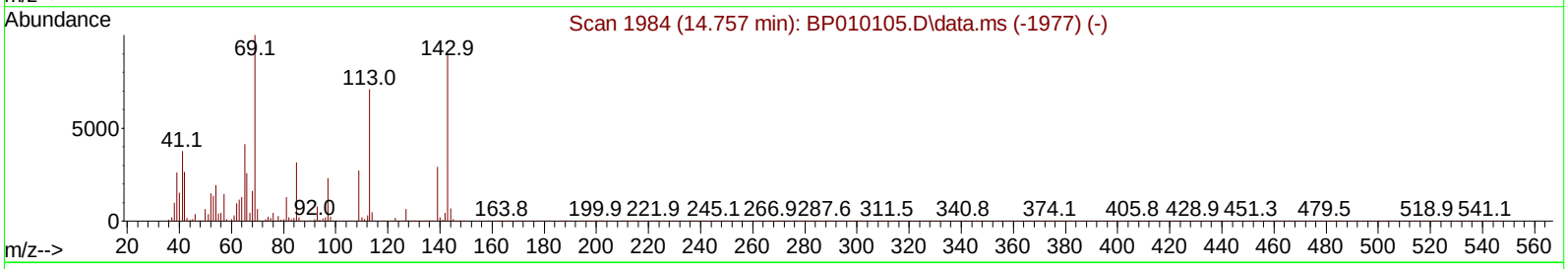

TIC: BP010111.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 3.45 ng/ul m

response 18403

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.77#
41.00	23.20	50.57#
42.00	18.00	31.37#

Quantitation Report (QT Reviewed)

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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	7.934	152	132213	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	554160	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	371965	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	791122	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	670545	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	506852	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	16006	5.279	ng/uL	0.00
4) Pyridine-d5	3.787	84	65267	7.652	ng/ul	0.00
7) Phenol-d5	7.093	99	67302	5.596	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.257	67	221801	30.737	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	192790	21.646	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	132990	13.347	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	138913m	31.518	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	133541	27.743	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	223412	23.841	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	291286	21.679	ng/ul	-0.01
46) Dimethylphthalate-d6	13.975	166	993116	34.323	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1058197	30.218	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	18403m	3.455	ng/ul	0.00
60) Fluorene-d10	15.557	176	834925	33.545	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	135971	27.181	ng/ul	0.00
73) Anthracene-d10	17.416	188	1393032	36.397	ng/ul	0.00
81) Pyrene-d10	19.645	212	1661188	43.562	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1052592	39.613	ng/ul	0.00
Target Compounds						Qvalue

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

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