

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110920\
 Data File : BP003894.D
 Acq On : 10 Nov 2020 08:41
 Operator : CG/JU
 Sample : L4617-05
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PE-5-(0.5-1.0)

Quant Time: Nov 10 09:43:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 06 11:33:14 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.305	152	195325	20.00	ng	0.00
21) Naphthalene-d8	11.134	136	768819	20.00	ng	# 0.00
39) Acenaphthene-d10	14.922	164	471764	20.00	ng	0.00
64) Phenanthrene-d10	17.651	188	888571	20.00	ng	# 0.00
76) Chrysene-d12	21.674	240	721804	20.00	ng	0.00
86) Perylene-d12	24.157	264	709498	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.828	112	1013761	86.62	ng	0.00
7) Phenol-d6	7.475	99	1308532	77.89	ng	0.00
23) Nitrobenzene-d5	9.469	82	833961	53.12	ng	0.00
42) 2,4,6-Tribromophenol	16.404	330	496518	84.15	ng	0.00
45) 2-Fluorobiphenyl	13.546	172	1728504	51.22	ng	0.00
79) Terphenyl-d14	20.145	244	1765711	47.26	ng	0.01
Target Compounds						
10) Phenol	7.499	94	32669	2.02	ng	Qvalue 69
50) Dimethylphthalate	14.345	163	296505	8.04	ng	98
78) Pyrene	19.969	202	106545	2.18	ng	97

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

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