

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102220\
 Data File : BP003715.D
 Acq On : 23 Oct 2020 09:12
 Operator : CG/JU
 Sample : L4484-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 72-11998

Quant Time: Oct 23 09:48:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 19 14:54:24 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.528	152	168316	20.00	ng	0.00
21) Naphthalene-d8	11.369	136	631921	20.00	ng	# 0.00
39) Acenaphthene-d10	15.122	164	426741	20.00	ng	0.00
64) Phenanthrene-d10	17.839	188	998333	20.00	ng	0.00
76) Chrysene-d12	21.851	240	1074314	20.00	ng	# 0.00
86) Perylene-d12	24.427	264	983301	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.028	112	1225119	121.71	ng	0.00
7) Phenol-d6	7.675	99	1423879	100.12	ng	0.00
23) Nitrobenzene-d5	9.693	82	1270400	85.05	ng	0.00
42) 2,4,6-Tribromophenol	16.598	330	916553	137.67	ng	0.00
45) 2-Fluorobiphenyl	13.757	172	2888456	87.50	ng	0.00
79) Terphenyl-d14	20.310	244	4949909	81.03	ng	0.00
Target Compounds						
10) Phenol	7.699	94	86493	6.39	ng	78
50) Dimethylphthalate	14.539	163	311207	8.88	ng	99

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

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