

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102320\
 Data File : BP003739.D
 Acq On : 24 Oct 2020 03:56
 Operator : CG/JU
 Sample : L4496-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SA-01-102220

Quant Time: Oct 24 05:40:56 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	108328	20.00	ng	0.00
21) Naphthalene-d8	11.369	136	382920	20.00	ng	# 0.00
39) Acenaphthene-d10	15.122	164	238502	20.00	ng	0.00
64) Phenanthrene-d10	17.845	188	529572	20.00	ng	0.00
76) Chrysene-d12	21.857	240	599582	20.00	ng	# 0.00
86) Perylene-d12	24.433	264	642717	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.028	112	822286	126.93	ng	0.00
7) Phenol-d6	7.675	99	923862	100.94	ng	0.00
23) Nitrobenzene-d5	9.693	82	837145	92.48	ng	0.00
42) 2,4,6-Tribromophenol	16.598	330	548059	147.29	ng	0.00
45) 2-Fluorobiphenyl	13.751	172	1806095	97.90	ng	0.00
79) Terphenyl-d14	20.310	244	3122089	91.58	ng	0.00
Target Compounds						
10) Phenol	7.705	94	48305	5.54	ng	Qvalue 79
50) Dimethylphthalate	14.540	163	178206	9.09	ng	100

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

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