

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101119\
 Data File : BE100635.D
 Acq On : 11 Oct 2019 10:50
 Operator : JU
 Sample : K4839-08DL 50X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 11 14:47:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 14:01:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	5317	0.40	ng	-0.01
7) Naphthalene-d8	10.64	136	20523	0.40	ng	0.00
13) Acenaphthene-d10	14.48	164	11975	0.40	ng	0.00
19) Phenanthrene-d10	17.20	188	28030	0.40	ng	0.00
27) Chrysene-d12	21.37	240	40151	0.40	ng	0.00
34) Perylene-d12	23.85	264	44341	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.44	112	49016	2.99	ng	0.00
5) Phenol-d6	7.01	99	74370	3.21	ng	0.00
8) Nitrobenzene-d5	8.99	82	28134	2.21	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	52	0.00	ng	0.00
14) 2,4,6-Tribromophenol	15.96	330	10090	3.19	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	102485	2.35	ng	0.00
25) Fluoranthene-d10	19.23	212	179	0.00	ng	0.00
29) Terphenyl-d14	19.83	244	202135	2.30	ng	0.00
Target Compounds						
22) Pentachlorophenol	16.85	266	8835	2.018	ng	99
32) Bis(2-ethylhexyl)phthalate	21.29	149	5634	0.021	ng	# 98

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

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