

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032318\
 Data File : BE095828.D
 Acq On : 23 Mar 2018 18:13
 Operator : SJ/JU
 Sample : PB107346BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB107346BL

Quant Time: Mar 24 00:19:42 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1205	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	4339	0.40	ng	0.00
13) Acenaphthene-d10	15.00	164	2323	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	5465	0.40	ng	0.00
27) Chrysene-d12	21.79	240	6438	0.40	ng	-0.01
34) Perylene-d12	24.43	264	6606	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.91	112	1434	0.36	ng	0.00
5) Phenol-d6	7.57	99	1876	0.34	ng	0.00
8) Nitrobenzene-d5	9.59	82	2119	0.41	ng	0.00
11) 2-Methylnaphthalene-d10	12.80	152	2264	0.31	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	385	0.29	ng	0.00
15) 2-Fluorobiphenyl	13.63	172	4267	0.44	ng	-0.01
25) Fluoranthene-d10	19.68	212	25387	0.31	ng	0.00
29) Terphenyl-d14	20.24	244	6207	0.45	ng	0.00
Target Compounds						
26) Fluoranthene	19.72	202	590	0.026	ng	# 95
28) Pyrene	20.06	202	821	0.031	ng	# 90
30) Benzo(a)anthracene	21.78	228	912	0.041	ng	# 81
31) Chrysene	21.84	228	827	0.041	ng	# 74
33) Indeno(1,2,3-cd)pyrene	27.24	276	1205	0.055	ng	# 92
39) Benzo(g,h,i)perylene	28.10	276	1067	0.055	ng	# 55

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

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