

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101918\
 Data File : BE097992.D
 Acq On : 19 Oct 2018 15:24
 Operator : SJ/JU
 Sample : J5237-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Oct 19 23:54:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	196456	0.40	ng	0.00
7) Naphthalene-d8	10.68	136	922052	0.40	ng	-0.01
13) Acenaphthene-d10	14.54	164	644434	0.40	ng	-0.01
19) Phenanthrene-d10	17.29	188	1517863	0.40	ng	-0.01
27) Chrysene-d12	21.48	240	1615844	0.40	ng	-0.01
34) Perylene-d12	24.04	264	1681895	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.45	112	763044	1.23	ng	-0.01
5) Phenol-d6	7.06	99	827388	0.90	ng	0.00
8) Nitrobenzene-d5	9.03	82	1494196	1.71	ng	-0.01
11) 2-Methylnaphthalene-d10	12.28	152	258	0.00	ng	-0.01
14) 2,4,6-Tribromophenol	16.04	330	909354	2.62	ng	0.00
15) 2-Fluorobiphenyl	13.18	172	3130536	1.19	ng	0.00
25) Fluoranthene-d10	19.39	212	756	0.00	ng	0.06
29) Terphenyl-d14	19.93	244	4683789	1.26	ng	0.00
Target Compounds						
22) Pentachlorophenol	16.95	266	1220	0.036	ng	Qvalue 96

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

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