

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092118\
 Data File : BE097617.D
 Acq On : 21 Sep 2018 18:32
 Operator : SJ/JU
 Sample : PB112764TB
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB112764TB

Manual Integrations
 APPROVED

Sohil
 9/24/2018 8:29:57 AM

Quant Time: Sep 22 04:35:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE091918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 10:54:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1006	0.40	ng	0.00
7) Naphthalene-d8	10.16	136	3920	0.40	ng	0.03
13) Acenaphthene-d10	14.03	164	2038	0.40	ng	0.02
19) Phenanthrene-d10	16.78	188	4991	0.40	ng	0.02
27) Chrysene-d12	20.98	240	6815	0.40	ng	0.01
34) Perylene-d12	23.22	264	6686	0.40	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.06	112	1155	0.38	ng	0.01
5) Phenol-d6	6.65	99	1326m	0.34	ng	0.05
8) Nitrobenzene-d5	8.59	82	868m	0.29	ng	0.06
11) 2-Methylnaphthalene-d10	11.76	152	2394	0.44	ng	0.05
14) 2,4,6-Tribromophenol	15.56	330	225m	0.32	ng	0.05
15) 2-Fluorobiphenyl	12.66	172	3255	0.45	ng	0.03
25) Fluoranthene-d10	18.82	212	27834	0.44	ng	0.01
29) Terphenyl-d14	19.42	244	4995	0.40	ng	0.00

Target Compounds Qvalue

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

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