

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100318\
 Data File : BE097717.D
 Acq On : 3 Oct 2018 18:40
 Operator : SJ/JU
 Sample : J4870-08DL 100X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 04 04:47:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 03 12:17:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	4799	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	20148	0.40	ng	0.00
13) Acenaphthene-d10	14.56	164	11397	0.40	ng	0.00
19) Phenanthrene-d10	17.30	188	31502	0.40	ng	0.00
27) Chrysene-d12	21.49	240	36274	0.40	ng	0.00
34) Perylene-d12	24.04	264	41058	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	16204	1.43	ng	0.00
5) Phenol-d6	7.05	99	24961	1.43	ng	0.00
8) Nitrobenzene-d5	9.04	82	7288	1.11	ng	0.00
11) 2-Methylnaphthalene-d10	12.32	152	22	0.00	ng	0.03
14) 2,4,6-Tribromophenol	16.04	330	3046	1.47	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	43388	0.86	ng	-0.02
25) Fluoranthene-d10	19.34	212	216	0.00	ng	0.00
29) Terphenyl-d14	19.93	244	77576	1.05	ng	0.00
Target Compounds						
9) Naphthalene	10.74	128	13236	0.065	ng	Qvalue # 97
22) Pentachlorophenol	16.95	266	3507	1.202	ng	96

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

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