

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101617\
 Data File : BE094423.D
 Acq On : 16 Oct 2017 19:01
 Operator : SJ/JU
 Sample : I5266-08DL 100X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-ACIDS-WPDL

Manual Integrations
 APPROVED

Sohil
 10/17/2017 5:14:35 PM

Quant Time: Oct 16 20:18:10 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 14:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1251	0.40	ng	-0.01
7) Naphthalene-d8	10.70	136	6729	0.40	ng	0.00
13) Acenaphthene-d10	14.53	164	3952	0.40	ng	0.00
19) Phenanthrene-d10	17.27	188	6749	0.40	ng	0.00
27) Chrysene-d12	21.43	240	9446	0.40	ng	0.00
34) Perylene-d12	23.96	264	8886	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.51	112	5003	1.18	ng	0.00
5) Phenol-d6	7.11	99	7813	1.18	ng	-0.01
8) Nitrobenzene-d5	9.06	82	5918	1.01	ng	-0.02
11) 2-Methylnaphthalene-d10	0.00	152	0d	0.00	ng	
14) 2,4,6-Tribromophenol	16.03	330	1727	1.15	ng	0.00
15) 2-Fluorobiphenyl	13.15	172	14060	1.00	ng	0.00
25) Fluoranthene-d10	0.00	212	0d	0.00	ng	
29) Terphenyl-d14	19.87	244	16927	0.89	ng	0.00
Target Compounds						Qvalue
22) Pentachlorophenol	16.97	266	837m	1.334	ng	

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

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