

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011017\
 Data File : BF092178.D
 Acq On : 10 Jan 2017 13:33
 Operator : UM/SJ
 Sample : I1045-03RE 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAP-SP1RE

Quant Time: Jan 10 15:44:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	162777	20.00	ng	-0.01
21) Naphthalene-d8	7.91	136	611005	20.00	ng	-0.02
38) Acenaphthene-d10	9.67	164	298522	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	419999	20.00	ng	-0.01
75) Chrysene-d12	13.79	240	276938	20.00	ng	0.00
86) Perylene-d12	15.17	264	194919	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	644394	66.10	ng	-0.01
7) Phenol-d6	6.30	99	662598	55.67	ng	-0.01
23) Nitrobenzene-d5	7.19	82	353485	32.17	ng	-0.02
41) 2,4,6-Tribromophenol	10.46	330	89105	36.07	ng	-0.01
44) 2-Fluorobiphenyl	9.00	172	632735	35.99	ng	-0.01
78) Terphenyl-d14	12.73	244	478553	41.91	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.31	94	37552	2.77	ng	77
49) Dimethylphthalate	9.40	163	98371	5.15	ng	98
70) Phenanthrene	11.17	178	104161	4.57	ng	97
74) Fluoranthene	12.36	202	134717	3.19	ng	97
77) Pyrene	12.59	202	129793	6.39	ng	98
80) Benzo(a)anthracene	13.77	228	46418m	2.97	ng	
82) Chrysene	13.81	228	57479m	3.67	ng	

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

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