

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109360.D
 Acq On : 27 Sep 2018 2:18
 Operator : JU/SJ
 Sample : J5053-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RFK-RMAB-MW08-GW

Quant Time: Sep 27 11:35:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	88224	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	287019	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	126372	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	246808	20.00	ng	0.00
75) Chrysene-d12	13.84	240	272047	20.00	ng	0.00
86) Perylene-d12	15.24	264	224391	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	102804	19.19	ng	0.04
7) Phenol-d6	6.37	99	334525	48.94	ng	0.01
23) Nitrobenzene-d5	7.28	82	398675	82.00	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	161378	129.54	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	704447	84.07	ng	0.00
78) Terphenyl-d14	12.80	244	788648	71.14	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	7.13	107	37683	6.476	ng	94
27) 2,4-Dimethylphenol	7.69	122	879716	230.596	ng	# 90
31) Naphthalene	8.03	128	4288932	319.776	ng	95
37) 2-Methylnaphthalene	8.71	142	709255	82.030	ng	97

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

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