

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088372.D
 Acq On : 25 Jun 2016 12:53
 Operator : UM/SJ
 Sample : H3602-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 08-B2(0-2)C

Quant Time: Jun 26 02:32:09 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	79878	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	317404	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	146808	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	253452	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	159826	20.00	ng	-0.06
86) Perylene-d12	15.05	264	152004	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	565593	121.05	ng	-0.02
7) Phenol-d6	6.23	99	808891	142.85	ng	-0.03
23) Nitrobenzene-d5	7.14	82	543070	99.91	ng	-0.03
41) 2,4,6-Tribromophenol	10.38	330	104052	67.12	ng	-0.06
44) 2-Fluorobiphenyl	8.92	172	828284	88.88	ng	-0.05
78) Terphenyl-d14	12.65	244	609881	86.83	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	9.32	163	124481	11.71	ng	99
70) Phenanthrene	11.10	178	49421	3.72	ng	99
74) Fluoranthene	12.27	202	58160	4.14	ng	97
77) Pyrene	12.50	202	47632	4.02	ng	99
80) Benzo(a)anthracene	13.68	228	22067	2.42	ng	# 96
82) Chrysene	13.73	228	18322m	2.14	ng	
87) Benzo(b)fluoranthene	14.70	252	24761m	2.34	ng	
89) Benzo(a)pyrene	15.01	252	17350	2.02	ng	# 89

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

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