

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071516\
 Data File : BF089039.D
 Acq On : 16 Jul 2016 4:04
 Operator : UM/SJ
 Sample : H4020-04
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-COMP

Quant Time: Jul 16 04:51:47 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 11 18:13:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	58012	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	229411	20.00	ng	-0.03
38) Acenaphthene-d10	9.71	164	89396	20.00	ng	-0.03
63) Phenanthrene-d10	11.19	188	133509	20.00	ng	-0.03
75) Chrysene-d12	13.81	240	111084	20.00	ng	-0.05
86) Perylene-d12	15.17	264	78447	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	519400	134.18	ng	-0.01
7) Phenol-d6	6.33	99	706255	141.21	ng	-0.02
23) Nitrobenzene-d5	7.24	82	376905	86.10	ng	-0.03
41) 2,4,6-Tribromophenol	10.50	330	100091	120.57	ng	-0.03
44) 2-Fluorobiphenyl	9.04	172	579417	102.38	ng	-0.03
78) Terphenyl-d14	12.76	244	339946	68.16	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.44	163	230348	36.17	ng	99
74) Fluoranthene	12.39	202	32319	4.37	ng	96
77) Pyrene	12.62	202	28285	3.00	ng	97
80) Benzo(a)anthracene	13.79	228	15728m	2.20	ng	
82) Chrysene	13.83	228	16856m	2.38	ng	
87) Benzo(b)fluoranthene	14.80	252	21614m	4.39	ng	
89) Benzo(a)pyrene	15.11	252	11900	2.86	ng	93

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

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