

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082316\
 Data File : BF089853.D
 Acq On : 24 Aug 2016 1:36
 Operator : UM/SJ
 Sample : H4547-09DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-7(3-3.5)DL

Quant Time: Aug 25 14:14:51 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 19 14:02:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	428861	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	1574687	20.00	ng	-0.02
38) Acenaphthene-d10	9.71	164	802431	20.00	ng	0.00
63) Phenanthrene-d10	11.19	188	1601413	20.00	ng	0.00
75) Chrysene-d12	13.84	240	1300041	20.00	ng	-0.03
86) Perylene-d12	15.33	264	1071377	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	513364	20.53	ng	-0.02
7) Phenol-d6	6.30	99	670653	21.85	ng	-0.02
23) Nitrobenzene-d5	7.23	82	410532	16.20	ng	-0.02
41) 2,4,6-Tribromophenol	10.49	330	119818	15.71	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	802437	17.58	ng	-0.01
78) Terphenyl-d14	12.76	244	714318	12.43	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.98	128	2374789	32.27	ng	99
37) 2-Methylnaphthalene	8.67	142	2468259	51.84	ng	97
49) Dimethylphthalate	9.44	163	216568	3.83	ng	# 94
51) Acenaphthene	9.74	154	119461	2.45	ng	# 86
57) Fluorene	10.25	166	319137	6.53	ng	# 82
70) Phenanthrene	11.21	178	603920	7.47	ng	95

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

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