

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082316\
 Data File : BF089851.D
 Acq On : 24 Aug 2016 00:42
 Operator : UM/SJ
 Sample : H4547-02DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D2-(3-3.5)DL

Quant Time: Aug 25 14:10:06 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	410016	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	1603271	20.00	ng	-0.02
38) Acenaphthene-d10	9.70	164	780403	20.00	ng	-0.01
63) Phenanthrene-d10	11.18	188	1580315	20.00	ng	-0.01
75) Chrysene-d12	13.82	240	1310150	20.00	ng	-0.06
86) Perylene-d12	15.27	264	1052367	20.00	ng	-0.10
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	675854	28.27	ng	-0.02
7) Phenol-d6	6.30	99	953916	32.50	ng	-0.02
23) Nitrobenzene-d5	7.23	82	502921	19.49	ng	-0.02
41) 2,4,6-Tribromophenol	10.49	330	198043	26.70	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	1040135	23.42	ng	-0.01
78) Terphenyl-d14	12.77	244	973719	16.81	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.98	128	2209543	29.49	ng	99
37) 2-Methylnaphthalene	8.67	142	2016938	41.61	ng	99
45) 1,1'-Biphenyl	9.13	154	147461	2.44	ng	95
49) Dimethylphthalate	9.43	163	268647	4.89	ng	# 97
57) Fluorene	10.25	166	122426	2.58	ng	# 78
70) Phenanthrene	11.20	178	209014	2.62	ng	94

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

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