

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\
 Data File : BF084918.D
 Acq On : 6 Feb 2016 18:12
 Operator : SJ/IZ
 Sample : H1338-04
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AW-1

Quant Time: Feb 08 02:58:26 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 14:16:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	175496	20.00	ng	-0.02
21) Naphthalene-d8	7.95	136	716984	20.00	ng	-0.02
38) Acenaphthene-d10	9.70	164	302024	20.00	ng	-0.03
63) Phenanthrene-d10	11.17	188	447039	20.00	ng	-0.03
75) Chrysene-d12	13.81	240	324160	20.00	ng	-0.02
86) Perylene-d12	15.17	264	313054	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	714842	63.45	ng	-0.02
7) Phenol-d6	6.34	99	577123	41.32	ng	0.00
23) Nitrobenzene-d5	7.23	82	1047982	82.34	ng	-0.03
41) 2,4,6-Tribromophenol	10.49	330	333169	105.76	ng	-0.03
44) 2-Fluorobiphenyl	9.04	172	1563140	88.31	ng	-0.02
78) Terphenyl-d14	12.76	244	1002970	70.11	ng	-0.03
Target Compounds						Qvalue
20) 3+4-Methylphenols	7.10	107	76306	6.10	ng	89
31) Naphthalene	7.97	128	564982	15.71	ng	99
37) 2-Methylnaphthalene	8.67	142	534668	23.75	ng	96
49) Dimethylphthalate	9.44	163	133185	5.79	ng	# 90
51) Acenaphthene	9.73	154	39558	2.02	ng	97
70) Phenanthrene	11.20	178	69229	2.79	ng	96

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

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