

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
 Data File : BF087689.D
 Acq On : 5 Jun 2016 5:10
 Operator : UM/SJ
 Sample : H3344-01
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-8

Manual Integrations
 APPROVED

SOHIL
 6/6/2016 7:18:50 PM

Quant Time: Jun 05 06:37:38 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 04 11:07:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	133556	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	315973	20.00	ng	0.01
38) Acenaphthene-d10	9.90	164	246378	20.00	ng	-0.01
63) Phenanthrene-d10	11.38	188	483878	20.00	ng	0.00
75) Chrysene-d12	14.01	240	384210	20.00	ng	-0.01
86) Perylene-d12	15.44	264	291469	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	506368	61.28	ng	0.00
7) Phenol-d6	6.48	99	415070	40.05	ng	-0.01
23) Nitrobenzene-d5	7.43	82	867823	158.90	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	327366	132.37	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1473955	95.78	ng	-0.01
78) Terphenyl-d14	12.97	244	1144479	72.69	ng	0.00
Target Compounds						
31) Naphthalene	8.22	128	8925942	581.01	ng	Qvalue 97
37) 2-Methylnaphthalene	8.86	142	829131	81.73	ng	# 97
45) 1,1'-Biphenyl	9.32	154	138046	6.90	ng	95
49) Dimethylphthalate	9.62	163	112092	6.26	ng	97
51) Acenaphthene	9.93	154	298259	18.68	ng	99
57) Fluorene	10.45	166	70618m	4.15	ng	
70) Phenanthrene	11.41	178	80939	3.09	ng	98

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

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