

Data Path : Z:\HPCHEM1\BNA F\DATA\BF053116\
 Data File : BF087598.D
 Acq On : 1 Jun 2016 00:52
 Operator : UM/SJ
 Sample : H3346-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-201605A

Quant Time: Jun 01 04:06:41 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 01 03:17:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	99233	20.00	ng	0.00
21) Naphthalene-d8	9.47	136	376641	20.00	ng	0.00
38) Acenaphthene-d10	12.30	164	194014	20.00	ng	-0.01
63) Phenanthrene-d10	14.71	188	400891	20.00	ng	0.00
75) Chrysene-d12	18.43	240	358229	20.00	ng	0.01
86) Perylene-d12	20.13	264	290112	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	766651	135.75	ng	0.01
7) Phenol-d6	7.00	99	926113	121.36	ng	-0.01
23) Nitrobenzene-d5	8.36	82	754137	100.91	ng	-0.01
41) 2,4,6-Tribromophenol	13.60	330	208679	111.93	ng	-0.01
44) 2-Fluorobiphenyl	11.24	172	1287038	93.52	ng	-0.01
78) Terphenyl-d14	17.06	244	1425408	84.59	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	9.54	128	10356429	548.74	ng	95
37) 2-Methylnaphthalene	10.64	142	801406	67.97	ng	96
45) 1,1'-Biphenyl	11.43	154	107430	6.58	ng	99
48) Acenaphthylene	12.07	152	744487	37.63	ng	98
57) Fluorene	13.22	166	103300	7.24	ng	98
70) Phenanthrene	14.77	178	171588	7.73	ng	97

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

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