

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010816\
 Data File : BF084346.D
 Acq On : 9 Jan 2016 1:26
 Operator : SJ/UM
 Sample : H1036-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-DRUM1

Quant Time: Jan 09 04:49:36 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 04 12:22:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	262502	20.00	ng	-0.05
21) Naphthalene-d8	8.14	136	1163084	20.00	ng	-0.05
38) Acenaphthene-d10	9.89	164	537076	20.00	ng	-0.06
63) Phenanthrene-d10	11.38	188	1016622	20.00	ng	-0.05
75) Chrysene-d12	14.01	240	729308	20.00	ng	-0.06
86) Perylene-d12	15.44	264	535050	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1924474	118.58	ng	-0.01
7) Phenol-d6	6.51	99	2013585	98.79	ng	-0.02
23) Nitrobenzene-d5	7.42	82	1715862	82.11	ng	-0.05
41) 2,4,6-Tribromophenol	10.69	330	760295	140.22	ng	-0.05
44) 2-Fluorobiphenyl	9.22	172	2547914	83.33	ng	-0.05
78) Terphenyl-d14	12.97	244	2709350	82.92	ng	-0.05
Target Compounds						Qvalue
31) Naphthalene	8.17	128	2514348	47.65	ng	98
32) Benzoic acid	7.88	122	9841	6.93	ng	# 82
37) 2-Methylnaphthalene	8.85	142	351325	9.27	ng	99
51) Acenaphthene	9.93	154	305127	9.18	ng	98
54) Dibenzofuran	10.10	168	282585	3.85	ng	99
57) Fluorene	10.44	166	204734	3.98	ng	98
70) Phenanthrene	11.40	178	318401	5.99	ng	99
76) Benzidine	12.72	184	66772	2.55	ng	98

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

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