

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031116\
 Data File : BF085399.D
 Acq On : 12 Mar 2016 5:16
 Operator : UM/SJ
 Sample : H1754-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W170TH-SB-21(0.5-3.0)

Manual Integrations
 APPROVED

UMANGI
 3/14/2016 10:45:56 AM

Quant Time: Mar 12 06:48:31 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	173485	20.00	ng	-0.05
21) Naphthalene-d8	8.22	136	648885	20.00	ng	-0.05
38) Acenaphthene-d10	9.98	164	295260	20.00	ng	-0.05
63) Phenanthrene-d10	11.46	188	474935	20.00	ng	-0.05
75) Chrysene-d12	14.11	240	316897	20.00	ng	-0.05
86) Perylene-d12	15.57	264	290831	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1331974	128.79	ng	-0.03
7) Phenol-d6	6.57	99	2025856	149.51	ng	-0.03
23) Nitrobenzene-d5	7.50	82	1112907	91.78	ng	-0.05
41) 2,4,6-Tribromophenol	10.77	330	348958	138.12	ng	-0.05
44) 2-Fluorobiphenyl	9.30	172	1785836	91.98	ng	-0.05
78) Terphenyl-d14	13.05	244	1119903	82.04	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.69	163	250852	11.07	ng	98
70) Phenanthrene	11.49	178	62668	2.52	ng	96
74) Fluoranthene	12.68	202	112085	4.49	ng	97
77) Pyrene	12.90	202	113693	4.77	ng	97
80) Benzo(a)anthracene	14.09	228	61741	3.25	ng	99
82) Chrysene	14.13	228	41466	2.37	ng	99
87) Benzo(b)fluoranthene	15.15	252	61640m	3.18	ng	
89) Benzo(a)pyrene	15.50	252	43343	2.70	ng	# 91

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

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