

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030816\
 Data File : BF085273.D
 Acq On : 8 Mar 2016 23:31
 Operator : UM/SJ
 Sample : H1684-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 48-BRIDGE-GRID-1

Manual Integrations
 APPROVED

UMANGI
 3/9/2016 9:30:52 AM

Quant Time: Mar 09 01:21:11 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	6.96	152	156071	20.00	ng	-0.02
21) Naphthalene-d8	8.24	136	593858	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	291720	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	541773	20.00	ng	-0.02
75) Chrysene-d12	14.12	240	390037	20.00	ng	-0.03
86) Perylene-d12	15.58	264	269688	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	1265470	136.01	ng	0.00
7) Phenol-d6	6.58	99	1259655	103.33	ng	-0.02
23) Nitrobenzene-d5	7.52	82	950548	85.65	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	379084	151.87	ng	-0.02
44) 2-Fluorobiphenyl	9.33	172	1824363	95.11	ng	-0.02
78) Terphenyl-d14	13.08	244	1839125	109.46	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.26	128	1078358	36.93	ng	99
37) 2-Methylnaphthalene	8.95	142	99715	5.32	ng	96
51) Acenaphthene	10.04	154	56468	3.15	ng	97
57) Fluorene	10.55	166	38370	2.08	ng	99
70) Phenanthrene	11.51	178	78812	2.78	ng	96

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

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