

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031116\
 Data File : BF085390.D
 Acq On : 11 Mar 2016 23:49
 Operator : UM/SJ
 Sample : H1743-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MNR-FT-ERM33

Manual Integrations
 APPROVED

UMANGI
 3/14/2016 10:45:43 AM

Quant Time: Mar 14 03:57:49 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	143866	20.00	ng	-0.05
21) Naphthalene-d8	8.22	136	550097	20.00	ng	-0.05
38) Acenaphthene-d10	9.98	164	249965	20.00	ng	-0.05
63) Phenanthrene-d10	11.46	188	403357	20.00	ng	-0.05
75) Chrysene-d12	14.11	240	275410	20.00	ng	-0.05
86) Perylene-d12	15.57	264	263844	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	583454	68.03	ng	-0.05
7) Phenol-d6	6.55	99	424173	37.75	ng	-0.06
23) Nitrobenzene-d5	7.50	82	877794	85.39	ng	-0.05
41) 2,4,6-Tribromophenol	10.77	330	276817	129.42	ng	-0.05
44) 2-Fluorobiphenyl	9.30	172	1549846	94.30	ng	-0.05
78) Terphenyl-d14	13.05	244	1020892	86.05	ng	-0.03
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.94	142	241008	13.88	ng	98
49) Dimethylphthalate	9.69	163	182789	9.53	ng	99
57) Fluorene	10.53	166	36543	2.31	ng	100
72) Carbazole	11.69	167	49669	2.52	ng	99

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

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