

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086410.D
 Acq On : 24 Apr 2016 00:42
 Operator : UM/SJ
 Sample : H2652-13 4X
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B6(0-6.5)-C

Manual Integrations
 APPROVED

sohil
 4/25/2016 12:36:15 PM

Quant Time: Apr 25 07:20:36 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 23 00:24:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	157063	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	606692	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	247474	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	437889	20.00	ng	0.00
75) Chrysene-d12	14.00	240	324207	20.00	ng	0.00
86) Perylene-d12	15.43	264	227387	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	266064	29.25	ng	0.01
7) Phenol-d6	6.48	99	346049	27.99	ng	0.00
23) Nitrobenzene-d5	7.40	82	176230	17.21	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	51696	23.08	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	335355	23.12	ng	0.00
78) Terphenyl-d14	12.94	244	287903	22.06	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.60	163	36362	2.03	ng	Qvalue 100
74) Fluoranthene	12.57	202	70759	2.98	ng	96
77) Pyrene	12.80	202	64905	2.87	ng	97
80) Benzo(a)anthracene	13.99	228	34683	2.06	ng	92
87) Benzo(b)fluoranthene	15.01	252	27111m	2.17	ng	

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

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