

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
 Data File : BF086648.D
 Acq On : 3 May 2016 16:57
 Operator : UM/SJ
 Sample : H2685-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2B-CS-6(10-20FTBGS)

Manual Integrations
 APPROVED

sohil
 5/4/2016 6:52:51 PM

Quant Time: May 04 04:03:34 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	193205	20.00	ng	-0.06
21) Naphthalene-d8	8.04	136	775219	20.00	ng	-0.06
38) Acenaphthene-d10	9.79	164	348638	20.00	ng	-0.06
63) Phenanthrene-d10	11.26	188	537048	20.00	ng	-0.06
75) Chrysene-d12	13.91	240	387938	20.00	ng	-0.06
86) Perylene-d12	15.30	264	373206	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1324241	118.22	ng	-0.04
7) Phenol-d6	6.40	99	1587403	101.54	ng	-0.05
23) Nitrobenzene-d5	7.32	82	1169542	81.32	ng	-0.06
41) 2,4,6-Tribromophenol	10.58	330	323703	88.04	ng	-0.06
44) 2-Fluorobiphenyl	9.12	172	1418108	67.94	ng	-0.06
78) Terphenyl-d14	12.85	244	908496	51.59	ng	-0.06
Target Compounds						Qvalue
49) Dimethylphthalate	9.52	163	358523	13.66	ng	99
70) Phenanthrene	11.29	178	121536	4.31	ng	99
74) Fluoranthene	12.48	202	259066	9.11	ng	96
77) Pyrene	12.71	202	234292	8.38	ng	99
80) Benzo(a)anthracene	13.89	228	129840	6.28	ng	99
82) Chrysene	13.93	228	83733m	3.72	ng	
87) Benzo(b)fluoranthene	14.90	252	138399m	6.30	ng	
89) Benzo(a)pyrene	15.24	252	87652	4.23	ng	97
91) Benzo(g,h,i)perylene	17.01	276	42076	2.48	ng	94

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

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