

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042916\
 Data File : BF086492.D
 Acq On : 29 Apr 2016 15:13
 Operator : UM/SJ
 Sample : H2685-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2A-CS-10(6-12FTBGS)

Manual Integrations
 APPROVED

sohil
 4/30/2016 11:22:45 AM

Quant Time: May 02 11:14:32 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	6.77	152	159097	20.00	ng	-0.03
21) Naphthalene-d8	8.06	136	681501	20.00	ng	-0.03
38) Acenaphthene-d10	9.81	164	324738	20.00	ng	-0.03
63) Phenanthrene-d10	11.30	188	585928	20.00	ng	-0.02
75) Chrysene-d12	13.93	240	402611	20.00	ng	-0.03
86) Perylene-d12	15.32	264	330459	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1234373	133.83	ng	-0.01
7) Phenol-d6	6.42	99	1629255	126.56	ng	-0.02
23) Nitrobenzene-d5	7.34	82	1195438	94.55	ng	-0.03
41) 2,4,6-Tribromophenol	10.60	330	414537	121.05	ng	-0.03
44) 2-Fluorobiphenyl	9.14	172	1804768	101.17	ng	-0.03
78) Terphenyl-d14	12.89	244	1572849	86.07	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.54	163	384454	15.73	ng	98
70) Phenanthrene	11.32	178	220933	7.19	ng	98
74) Fluoranthene	12.50	202	254558	8.20	ng	97
77) Pyrene	12.73	202	213921	7.37	ng	99
80) Benzo(a)anthracene	13.92	228	76483	3.56	ng	98
82) Chrysene	13.95	228	74285m	3.18	ng	
87) Benzo(b)fluoranthene	14.92	252	74791m	3.85	ng	
89) Benzo(a)pyrene	15.27	252	52174	2.84	ng	96

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

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