

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
 Data File : BF086645.D
 Acq On : 3 May 2016 15:22
 Operator : UM/SJ
 Sample : H2685-03
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2A-CS-2(0-6FTBGS)

Manual Integrations
 APPROVED

sohil
 5/4/2016 6:52:27 PM

Quant Time: May 04 03:13:23 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	191204	20.00	ng	-0.06
21) Naphthalene-d8	8.04	136	769235	20.00	ng	-0.06
38) Acenaphthene-d10	9.79	164	369058	20.00	ng	-0.06
63) Phenanthrene-d10	11.26	188	568176	20.00	ng	-0.06
75) Chrysene-d12	13.91	240	380103	20.00	ng	-0.06
86) Perylene-d12	15.30	264	377826	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1156368	104.32	ng	-0.04
7) Phenol-d6	6.40	99	1661245	107.38	ng	-0.05
23) Nitrobenzene-d5	7.32	82	1177458	82.51	ng	-0.06
41) 2,4,6-Tribromophenol	10.58	330	364429	93.64	ng	-0.06
44) 2-Fluorobiphenyl	9.12	172	1570355	71.90	ng	-0.06
78) Terphenyl-d14	12.85	244	1141624	66.17	ng	-0.06
Target Compounds						Qvalue
49) Dimethylphthalate	9.52	163	369434	13.30	ng	99
70) Phenanthrene	11.29	178	106420	3.57	ng	99
74) Fluoranthene	12.48	202	223447	7.42	ng	94
77) Pyrene	12.71	202	208568	7.62	ng	98
80) Benzo(a)anthracene	13.89	228	86624	4.27	ng	97
82) Chrysene	13.93	228	92662m	4.21	ng	
85) Indeno(1,2,3-cd)pyrene	16.63	276	57647	3.53	ng	93
87) Benzo(b)fluoranthene	14.91	252	139243m	6.26	ng	
89) Benzo(a)pyrene	15.24	252	90099	4.29	ng	# 95
91) Benzo(g,h,i)perylene	17.01	276	54503	3.17	ng	98

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

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