

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092916\
 Data File : BF090291.D
 Acq On : 29 Sep 2016 14:20
 Operator : UM/SJ
 Sample : H4982-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-17

Manual Integrations
 APPROVED

sohil
 9/30/2016 4:00:55 PM

Quant Time: Sep 29 15:51:16 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 26 15:58:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.46	152	136477	20.00	ng	-0.02
21) Naphthalene-d8	7.75	136	525415	20.00	ng	-0.02
38) Acenaphthene-d10	9.50	164	278745	20.00	ng	-0.01
63) Phenanthrene-d10	10.97	188	358698	20.00	ng	-0.01
75) Chrysene-d12	13.58	240	262361	20.00	ng	-0.02
86) Perylene-d12	14.90	264	221567	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.03	112	1167830	139.48	ng	0.00
7) Phenol-d6	6.12	99	1409877	136.35	ng	-0.01
23) Nitrobenzene-d5	7.04	82	981467	108.28	ng	-0.02
41) 2,4,6-Tribromophenol	10.28	330	287324	120.15	ng	-0.01
44) 2-Fluorobiphenyl	8.83	172	1283244	90.38	ng	-0.02
78) Terphenyl-d14	12.56	244	935292	90.51	ng	0.00
Target Compounds						
10) Phenol	6.14	94	30356	2.48	ng	Qvalue 85
49) Dimethylphthalate	9.23	163	374034	21.07	ng	99
70) Phenanthrene	10.98	178	55542	3.31	ng	99
74) Fluoranthene	12.16	202	99858	5.55	ng	95
77) Pyrene	12.39	202	105261	5.86	ng	97
80) Benzo(a)anthracene	13.57	228	48602	3.44	ng	# 87
82) Chrysene	13.60	228	42777m	3.11	ng	
87) Benzo(b)fluoranthene	14.57	252	60459m	4.93	ng	
89) Benzo(a)pyrene	14.86	252	34852	3.02	ng	# 85
91) Benzo(g,h,i)perylene	16.39	276	21732	2.47	ng	# 89

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

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