

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042916\
 Data File : BF086502.D
 Acq On : 29 Apr 2016 19:58
 Operator : UM/SJ
 Sample : H2757-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B1(2-4)

Manual Integrations
 APPROVED

sohil
 4/30/2016 11:22:49 AM

Quant Time: Apr 30 01:37:10 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	152667	20.00	ng	-0.03
21) Naphthalene-d8	8.06	136	661797	20.00	ng	-0.03
38) Acenaphthene-d10	9.81	164	275855	20.00	ng	-0.03
63) Phenanthrene-d10	11.30	188	426635	20.00	ng	-0.02
75) Chrysene-d12	13.93	240	334033	20.00	ng	-0.03
86) Perylene-d12	15.32	264	268395	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	92844	10.49	ng	-0.03
7) Phenol-d6	6.41	99	137739	11.15	ng	-0.03
23) Nitrobenzene-d5	7.33	82	77702	6.33	ng	-0.05
41) 2,4,6-Tribromophenol	10.60	330	22133	7.61	ng	-0.03
44) 2-Fluorobiphenyl	9.14	172	176110	3.79	ng	-0.03
78) Terphenyl-d14	12.88	244	108610	7.16	ng	-0.03
Target Compounds						Qvalue
70) Phenanthrene	11.32	178	75600	3.38	ng	99
74) Fluoranthene	12.50	202	119680	5.30	ng	98
77) Pyrene	12.73	202	112761	4.68	ng	98
80) Benzo(a)anthracene	13.92	228	66527	3.74	ng	96
82) Chrysene	13.95	228	46933m	2.42	ng	
87) Benzo(b)fluoranthene	14.93	252	57405m	3.64	ng	
89) Benzo(a)pyrene	15.27	252	43415	2.91	ng	# 87

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

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