

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092316\
 Data File : BF090204.D
 Acq On : 23 Sep 2016 18:16
 Operator : UM/SJ
 Sample : H4867-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-WSG-16121-07

Manual Integrations
 APPROVED

sohil
 9/26/2016 5:58:11 PM

Quant Time: Sep 23 19:02:25 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 23 18:09:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	121776	20.00	ng	-0.03
21) Naphthalene-d8	7.78	136	452478	20.00	ng	-0.03
38) Acenaphthene-d10	9.52	164	222498	20.00	ng	-0.03
63) Phenanthrene-d10	10.98	188	261545	20.00	ng	-0.05
75) Chrysene-d12	13.60	240	237200	20.00	ng	-0.03
86) Perylene-d12	14.93	264	268241	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	960270	128.53	ng	-0.01
7) Phenol-d6	6.16	99	1255844	136.11	ng	-0.01
23) Nitrobenzene-d5	7.06	82	693767	88.88	ng	-0.03
41) 2,4,6-Tribromophenol	10.31	330	231331	121.19	ng	-0.03
44) 2-Fluorobiphenyl	8.86	172	930046	82.06	ng	-0.03
78) Terphenyl-d14	12.57	244	598042	64.01	ng	-0.03
Target Compounds						
22) Acetophenone	6.91	105	96006	8.80	ng	# 96
49) Dimethylphthalate	9.26	163	335046	23.64	ng	99
74) Fluoranthene	12.18	202	51184	3.90	ng	98
77) Pyrene	12.41	202	47187	2.91	ng	97
80) Benzo(a)anthracene	13.59	228	26898	2.11	ng	# 85
87) Benzo(b)fluoranthene	14.58	252	39417m	2.65	ng	

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

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