

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092316\
 Data File : BF090203.D
 Acq On : 23 Sep 2016 17:49
 Operator : UM/SJ
 Sample : H4867-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-WSG-16121-06

Manual Integrations
 APPROVED

sohil
 9/26/2016 5:58:14 PM

Quant Time: Sep 23 18:58:09 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	124079	20.00	ng	-0.03
21) Naphthalene-d8	7.78	136	463064	20.00	ng	-0.03
38) Acenaphthene-d10	9.52	164	237693	20.00	ng	-0.03
63) Phenanthrene-d10	10.98	188	300242	20.00	ng	-0.05
75) Chrysene-d12	13.60	240	238816	20.00	ng	-0.03
86) Perylene-d12	14.93	264	264012	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	998307	131.15	ng	-0.01
7) Phenol-d6	6.16	99	1389675	147.82	ng	-0.01
23) Nitrobenzene-d5	7.06	82	719137	90.03	ng	-0.03
41) 2,4,6-Tribromophenol	10.31	330	269620	132.22	ng	-0.03
44) 2-Fluorobiphenyl	8.86	172	982802	81.17	ng	-0.03
78) Terphenyl-d14	12.57	244	594220	63.17	ng	-0.03
Target Compounds						
22) Acetophenone	6.91	105	136533	12.23	ng	Qvalue # 95
49) Dimethylphthalate	9.26	163	405452	26.78	ng	98
74) Fluoranthene	12.18	202	58332	3.87	ng	96
77) Pyrene	12.41	202	49238	3.01	ng	97
80) Benzo(a)anthracene	13.59	228	35380m	2.75	ng	
82) Chrysene	13.62	228	40649m	3.25	ng	
85) Indeno(1,2,3-cd)pyrene	16.07	276	26427	2.61	ng	# 91
87) Benzo(b)fluoranthene	14.58	252	53935m	3.69	ng	
89) Benzo(a)pyrene	14.87	252	33875	2.46	ng	# 84
91) Benzo(g,h,i)perylene	16.42	276	23448	2.23	ng	# 86

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

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