

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092016\
 Data File : BF090218.D
 Acq On : 26 Sep 2016 13:51
 Operator : UM/SJ
 Sample : H4927-01DL 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-18ADL

Manual Integrations
 APPROVED

sohil
 9/27/2016 1:07:00 PM

Quant Time: Sep 26 16:10:05 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	6.48	152	153679	20.00	ng	0.00
21) Naphthalene-d8	7.77	136	621838	20.00	ng	0.00
38) Acenaphthene-d10	9.51	164	334622	20.00	ng	0.00
63) Phenanthrene-d10	10.98	188	528694	20.00	ng	0.00
75) Chrysene-d12	13.59	240	374390	20.00	ng	-0.01
86) Perylene-d12	14.91	264	318776	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.03	112	43066	4.57	ng	0.00
7) Phenol-d6	6.12	99	93766	8.05	ng	-0.01
23) Nitrobenzene-d5	7.05	82	54148	5.05	ng	-0.01
41) 2,4,6-Tribromophenol	10.30	330	4776	1.66	ng	0.00
44) 2-Fluorobiphenyl	8.84	172	142479	8.36	ng	-0.01
78) Terphenyl-d14	12.56	244	119833	8.13	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.24	163	43315	2.03	ng	98
51) Acenaphthene	9.54	154	96018	5.64	ng	98
57) Fluorene	10.04	166	80813	4.79	ng	100
70) Phenanthrene	11.01	178	965905	39.07	ng	99
71) Anthracene	11.05	178	300858	11.60	ng	98
74) Fluoranthene	12.18	202	565280	21.30	ng	92
77) Pyrene	12.40	202	923596	36.05	ng	# 96
80) Benzo(a)anthracene	13.59	228	420833	20.89	ng	98
82) Chrysene	13.62	228	374101m	19.07	ng	
85) Indeno(1,2,3-cd)pyrene	16.06	276	81604	5.14	ng	91
87) Benzo(b)fluoranthene	14.57	252	193825m	10.98	ng	
88) Benzo(k)fluoranthene	14.58	252	78423m	4.73	ng	
89) Benzo(a)pyrene	14.86	252	189481	11.41	ng	# 93
90) Dibenzo(a,h)anthracene	16.07	278	27931	2.16	ng	# 90
91) Benzo(a,h,i)perylene	16.39	276	93536	7.38	ng	96

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

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