

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088405.D
 Acq On : 26 Jun 2016 7:00
 Operator : UM/SJ
 Sample : H3598-10DL 10X
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PM-STA-1381(0-4)DL

Quant Time: Jun 26 08:26:52 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	88902	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	364535	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	160439	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	263242	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	197620	20.00	ng	-0.06
86) Perylene-d12	15.05	264	167776	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	33823	6.50	ng	-0.02
7) Phenol-d6	6.24	99	51951	8.24	ng	-0.02
23) Nitrobenzene-d5	7.14	82	31078	4.98	ng	-0.03
41) 2,4,6-Tribromophenol	10.38	330	8946	5.28	ng	-0.06
44) 2-Fluorobiphenyl	8.92	172	69419	1.35	ng	-0.05
78) Terphenyl-d14	12.64	244	47223	5.44	ng	-0.06
Target Compounds						Qvalue
48) Acenaphthylene	9.45	152	77194	4.67	ng	99
70) Phenanthrene	11.09	178	102382	7.41	ng	98
71) Anthracene	11.14	178	80607	5.79	ng	99
74) Fluoranthene	12.27	202	441843	30.31	ng	97
77) Pyrene	12.50	202	456549	31.13	ng	98
80) Benzo(a)anthracene	13.68	228	284300	25.25	ng	96
82) Chrysene	13.71	228	216551m	20.44	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	82448	8.38	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	252254m	21.57	ng	
88) Benzo(k)fluoranthene	14.70	252	127498m	14.45	ng	
89) Benzo(a)pyrene	14.99	252	163676	17.30	ng	97
90) Dibenzo(a,h)anthracene	16.27	278	21747m	2.47	ng	
91) Benzo(a,h,i)perylene	16.65	276	67062	7.42	ng	97

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

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