

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088401.D
 Acq On : 26 Jun 2016 5:05
 Operator : UM/SJ
 Sample : H3598-13DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PM-STA-1388(0-4)DL

Quant Time: Jun 26 06:31:03 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	74758m	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	292163	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	126638	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	197186	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	162396	20.00	ng	-0.06
86) Perylene-d12	15.04	264	127680	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.14	112	117133	26.79	ng	-0.03
7) Phenol-d6	6.23	99	142333	26.86	ng	-0.03
23) Nitrobenzene-d5	7.14	82	85966	17.18	ng	-0.03
41) 2,4,6-Tribromophenol	10.38	330	25745	19.25	ng	-0.06
44) 2-Fluorobiphenyl	8.92	172	181965	18.22	ng	-0.05
78) Terphenyl-d14	12.64	244	114639	16.06	ng	-0.06
Target Compounds						Qvalue
70) Phenanthrene	11.08	178	167896	16.23	ng	98
71) Anthracene	11.14	178	49296	4.72	ng	96
74) Fluoranthene	12.27	202	228629	20.94	ng	96
77) Pyrene	12.49	202	186965	15.51	ng	97
80) Benzo(a)anthracene	13.68	228	108379	11.71	ng	97
82) Chrysene	13.71	228	83356	9.57	ng	97
85) Indeno(1,2,3-cd)pyrene	16.29	276	31135	3.85	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	86141m	9.68	ng	
88) Benzo(k)fluoranthene	14.70	252	44819m	6.68	ng	
89) Benzo(a)pyrene	14.99	252	65442	9.09	ng	98
91) Benzo(g,h,i)perylene	16.65	276	28906	4.20	ng	95

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

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