

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\
 Data File : BF088460.D
 Acq On : 27 Jun 2016 19:27
 Operator : UM/SJ
 Sample : H3747-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C8-B4(0-4.5)C

Quant Time: Jun 28 06:13:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 19:14:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.54	152	64731m	20.00	ng	-0.01
21) Naphthalene-d8	7.83	136	254738	20.00	ng	-0.01
38) Acenaphthene-d10	9.56	164	106346	20.00	ng	-0.02
63) Phenanthrene-d10	11.05	188	182341	20.00	ng	-0.01
75) Chrysene-d12	13.68	240	118265	20.00	ng	-0.01
86) Perylene-d12	15.03	264	131516	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	415591	109.76	ng	-0.01
7) Phenol-d6	6.20	99	479870	104.57	ng	-0.03
23) Nitrobenzene-d5	7.12	82	289757	66.42	ng	-0.02
41) 2,4,6-Tribromophenol	10.36	330	96815	86.22	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	533639	78.40	ng	-0.02
78) Terphenyl-d14	12.63	244	357777	68.84	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.31	163	29609	3.84	ng	99
70) Phenanthrene	11.07	178	21362	2.23	ng	98
74) Fluoranthene	12.26	202	41829	4.14	ng	96
77) Pyrene	12.48	202	42847	4.88	ng	97
80) Benzo(a)anthracene	13.67	228	22072	3.28	ng	98
82) Chrysene	13.71	228	19557	3.08	ng	98
85) Indeno(1,2,3-cd)pyrene	16.30	276	15107	2.56	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	26173m	2.86	ng	
89) Benzo(a)pyrene	14.99	252	20380	2.75	ng	98
91) Benzo(g,h,i)perylene	16.66	276	15248	2.15	ng	97

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

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