

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\
 Data File : BF088481.D
 Acq On : 28 Jun 2016 7:02
 Operator : UM/SJ
 Sample : H3624-20DL 2X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1506(0-4)DL

Quant Time: Jun 28 07:41:57 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 28 05:25:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.54	152	75395m	20.00	ng	-0.01
21) Naphthalene-d8	7.83	136	313007	20.00	ng	-0.01
38) Acenaphthene-d10	9.58	164	127002	20.00	ng	-0.01
63) Phenanthrene-d10	11.05	188	192713	20.00	ng	-0.01
75) Chrysene-d12	13.69	240	158888	20.00	ng	0.00
86) Perylene-d12	15.04	264	168697	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	234897	53.26	ng	-0.01
7) Phenol-d6	6.20	99	309896	57.98	ng	-0.02
23) Nitrobenzene-d5	7.12	82	165534	30.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	43880	32.72	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	321320	36.59	ng	-0.02
78) Terphenyl-d14	12.64	244	193170	27.67	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	9.44	152	95650	7.32	ng	98
49) Dimethylphthalate	9.31	163	57618	6.26	ng	99
70) Phenanthrene	11.07	178	84944	8.40	ng	98
71) Anthracene	11.13	178	101279	9.93	ng	98
74) Fluoranthene	12.26	202	432249	40.50	ng	99
77) Pyrene	12.49	202	536326	45.49	ng	99
80) Benzo(a)anthracene	13.68	228	350346	38.69	ng	# 91
82) Chrysene	13.71	228	222214m	26.09	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	100095	12.65	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	328319m	27.92	ng	
88) Benzo(k)fluoranthene	14.70	252	132585m	14.95	ng	
89) Benzo(a)pyrene	14.99	252	216572	22.77	ng	97
90) Dibenzo(a,h)anthracene	16.27	278	24718	2.80	ng	# 75
91) Benzo(g,h,i)perylene	16.65	276	86662	9.54	ng	99

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

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