

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101250.D
 Acq On : 8 Dec 2017 21:46
 Operator : SJ/JU
 Sample : I6656-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BART-3-E(0-1)

Quant Time: Dec 08 23:36:15 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 08 18:17:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	46009	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	167684	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	69867	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	121030	20.00	ng	0.00
75) Chrysene-d12	14.00	240	100072	20.00	ng	0.00
86) Perylene-d12	15.49	264	75732	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	251888	90.47	ng	0.00
7) Phenol-d6	6.46	99	306990	90.81	ng	0.00
23) Nitrobenzene-d5	7.39	82	183399	65.24	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	59189	70.52	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	383545	75.11	ng	0.00
78) Terphenyl-d14	12.94	244	327232	71.52	ng	0.00
Target Compounds						
10) Phenol	6.47	94	8333	2.217	ng	91
49) Dimethylphthalate	9.57	163	33970	6.029	ng	100
70) Phenanthrene	11.38	178	56882	8.534	ng	98
74) Fluoranthene	12.57	202	104088	14.089	ng	96
77) Pyrene	12.80	202	85024	12.313	ng	100
80) Benzo(a)anthracene	13.99	228	35983m	5.695	ng	
82) Chrysene	14.03	228	40942	6.977	ng	98
83) Bis(2-ethylhexyl)phthalate	13.97	149	11619	2.677	ng	# 92
85) Indeno(1,2,3-cd)pyrene	16.97	276	14829	2.842	ng	# 93
87) Benzo(b)fluoranthene	15.05	252	39902m	8.796	ng	
88) Benzo(k)fluoranthene	15.07	252	12723m	2.847	ng	
89) Benzo(a)pyrene	15.42	252	24726	5.996	ng	# 85
91) Benzo(a,h,i)perylene	17.43	276	15365	4.484	ng	# 88

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

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