

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011017\
 Data File : BF092187.D
 Acq On : 10 Jan 2017 17:37
 Operator : UM/SJ
 Sample : I1057-07RE
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P3-SP9RE

Quant Time: Jan 11 12:13:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	159756	20.00	ng	-0.01
21) Naphthalene-d8	7.92	136	646159	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	290827	20.00	ng	0.00
63) Phenanthrene-d10	11.16	188	402258	20.00	ng	0.00
75) Chrysene-d12	13.80	240	239124	20.00	ng	0.01
86) Perylene-d12	15.19	264	159054	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	778326	81.35	ng	0.00
7) Phenol-d6	6.31	99	1126031	96.40	ng	0.00
23) Nitrobenzene-d5	7.20	82	556352	47.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.47	330	109303	45.41	ng	0.00
44) 2-Fluorobiphenyl	9.01	172	1066723	73.39	ng	0.00
78) Terphenyl-d14	12.74	244	643184	65.23	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.32	94	83083	6.24	ng	# 72
31) Naphthalene	7.94	128	110686	3.74	ng	99
49) Dimethylphthalate	9.41	163	69605	3.74	ng	99
51) Acenaphthene	9.70	154	105100	6.39	ng	99
54) Dibenzofuran	9.89	168	100874	4.54	ng	91
57) Fluorene	10.22	166	99115	6.14	ng	96
70) Phenanthrene	11.18	178	433925	19.89	ng	97
71) Anthracene	11.24	178	116070	4.12	ng	96
74) Fluoranthene	12.37	202	348047	15.70	ng	99
77) Pyrene	12.60	202	295550	16.85	ng	98
80) Benzo(a)anthracene	13.79	228	123763	9.17	ng	# 93
82) Chrysene	13.82	228	102513m	7.57	ng	
85) Indeno(1,2,3-cd)pyrene	16.48	276	37231	3.17	ng	98
87) Benzo(b)fluoranthene	14.80	252	95439m	9.64	ng	
88) Benzo(k)fluoranthene	14.81	252	47064m	5.57	ng	
89) Benzo(a)pyrene	15.13	252	63880	7.27	ng	# 59
91) Benzo(a,h,i)perylene	16.87	276	32263	4.29	ng	# 66

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

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