

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011017\
 Data File : BF092185.D
 Acq On : 10 Jan 2017 16:42
 Operator : UM/SJ
 Sample : I1057-05RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P3-SP7RE

Quant Time: Jan 11 12:10:46 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	144981	20.00	ng	-0.01
21) Naphthalene-d8	7.92	136	579072	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	256707	20.00	ng	0.00
63) Phenanthrene-d10	11.16	188	378994	20.00	ng	0.00
75) Chrysene-d12	13.80	240	226452	20.00	ng	0.01
86) Perylene-d12	15.19	264	161236	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	886550	102.10	ng	0.00
7) Phenol-d6	6.31	99	1301364	122.76	ng	0.00
23) Nitrobenzene-d5	7.20	82	699517	67.17	ng	-0.01
41) 2,4,6-Tribromophenol	10.47	330	127210	59.88	ng	0.00
44) 2-Fluorobiphenyl	9.01	172	1214690	99.09	ng	0.00
78) Terphenyl-d14	12.75	244	841626	90.14	ng	0.00
Target Compounds						
10) Phenol	6.32	94	87638	7.25	ng	Qvalue # 73
49) Dimethylphthalate	9.41	163	108873	6.63	ng	99
51) Acenaphthene	9.70	154	58959	4.06	ng	97
54) Dibenzofuran	9.88	168	52151	2.66	ng	# 95
57) Fluorene	10.22	166	54099	3.79	ng	# 93
70) Phenanthrene	11.18	178	298030	14.50	ng	98
71) Anthracene	11.24	178	66385	2.25	ng	97
74) Fluoranthene	12.37	202	305069	14.31	ng	98
77) Pyrene	12.60	202	277170	16.68	ng	98
80) Benzo(a)anthracene	13.79	228	114524	8.96	ng	# 91
82) Chrysene	13.82	228	87853m	6.85	ng	
85) Indeno(1,2,3-cd)pyrene	16.48	276	29182	2.63	ng	# 95
87) Benzo(b)fluoranthene	14.80	252	85593m	8.53	ng	
88) Benzo(k)fluoranthene	14.81	252	37542m	4.39	ng	
89) Benzo(a)pyrene	15.13	252	59574	6.69	ng	# 52
91) Benzo(g,h,i)perylene	16.88	276	29803	3.91	ng	# 53

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

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