

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\
 Data File : BF101167.D
 Acq On : 6 Dec 2017 15:04
 Operator : SJ/JU
 Sample : I6702-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 12/7/2017 11:35:38 AM

Quant Time: Dec 06 18:14:57 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 04 13:07:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	47074	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	185848	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	89171	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	178472	20.00	ng	0.00
75) Chrysene-d12	14.02	240	109311	20.00	ng	0.00
86) Perylene-d12	15.49	264	98879	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	316949	111.26	ng	0.00
7) Phenol-d6	6.48	99	376061	108.73	ng	0.00
23) Nitrobenzene-d5	7.41	82	237817	76.33	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	108455	101.25	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	511872	78.54	ng	0.00
78) Terphenyl-d14	12.96	244	479705	95.99	ng	0.00
Target Compounds						
10) Phenol	6.49	94	12006	3.122	ng	94
49) Dimethylphthalate	9.59	163	38318	5.328	ng	# 97
70) Phenanthrene	11.40	178	23919	2.434	ng	98
74) Fluoranthene	12.59	202	71025	6.519	ng	94
77) Pyrene	12.82	202	76259	10.110	ng	97
80) Benzo(a)anthracene	14.00	228	31326	4.539	ng	95
82) Chrysene	14.04	228	30511	4.760	ng	97
85) Indeno(1,2,3-cd)pyrene	16.97	276	22267	3.907	ng	# 90
87) Benzo(b)fluoranthene	15.06	252	34932m	5.898	ng	
89) Benzo(a)pyrene	15.43	252	25771	4.786	ng	98
91) Benzo(g,h,i)perylene	17.42	276	22341	4.994	ng	# 88

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

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