

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101787.D
 Acq On : 28 Dec 2017 3:37
 Operator : SJ/JU
 Sample : I7069-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-23-COMP

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:24 PM

Quant Time: Dec 28 06:43:28 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	167357	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	643008	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	253447	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	388021	20.00	ng	0.00
75) Chrysene-d12	13.97	240	311388	20.00	ng	0.00
86) Perylene-d12	15.45	264	282135	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1111827	98.96	ng	0.01
7) Phenol-d6	6.44	99	1329227	95.06	ng	0.00
23) Nitrobenzene-d5	7.36	82	827785	71.66	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	230034	94.19	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1153455	70.60	ng	0.00
78) Terphenyl-d14	12.90	244	743036	57.53	ng	0.00
Target Compounds						
10) Phenol	6.45	94	50197	3.150	ng	Qvalue # 69
49) Dimethylphthalate	9.55	163	139661	6.851	ng	98
70) Phenanthrene	11.34	178	393035	18.214	ng	99
71) Anthracene	11.40	178	105122	4.769	ng	98
74) Fluoranthene	12.54	202	589571	26.030	ng	97
77) Pyrene	12.77	202	549448	23.511	ng	100
80) Benzo(a)anthracene	13.97	228	311760	15.162	ng	97
82) Chrysene	14.00	228	301971	15.554	ng	98
85) Indeno(1,2,3-cd)pyrene	16.95	276	113335	6.556	ng	# 88
87) Benzo(b)fluoranthene	15.02	252	298454m	17.355	ng	
88) Benzo(k)fluoranthene	15.05	252	110620m	6.616	ng	
89) Benzo(a)pyrene	15.39	252	219746	13.862	ng	96
91) Benzo(a,h,i)perylene	17.40	276	96457	7.157	ng	# 92

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

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