

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101270.D
 Acq On : 9 Dec 2017 8:24
 Operator : SJ/JU
 Sample : I6744-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

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

Quant Time: Dec 11 14:38:29 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	44401	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	161971	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	66265	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	109487	20.00	ng	0.00
75) Chrysene-d12	14.01	240	88357	20.00	ng	0.00
86) Perylene-d12	15.49	264	66099	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	255757	95.18	ng	0.00
7) Phenol-d6	6.47	99	321234	98.47	ng	0.00
23) Nitrobenzene-d5	7.39	82	178727	65.82	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	55040	69.14	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	311866	64.39	ng	0.00
78) Terphenyl-d14	12.94	244	232311	57.51	ng	0.00
Target Compounds						
10) Phenol	6.47	94	11370	3.134	ng	Qvalue # 69
49) Dimethylphthalate	9.58	163	25084	4.694	ng	98
70) Phenanthrene	11.38	178	78710	13.054	ng	97
71) Anthracene	11.43	178	28707	4.704	ng	98
74) Fluoranthene	12.57	202	167459	25.056	ng	96
77) Pyrene	12.81	202	183693	30.129	ng	99
80) Benzo(a)anthracene	14.00	228	86535m	15.512	ng	
82) Chrysene	14.03	228	82707m	15.963	ng	
85) Indeno(1,2,3-cd)pyrene	16.97	276	28179	6.118	ng	# 90
87) Benzo(b)fluoranthene	15.05	252	75222m	19.000	ng	
88) Benzo(k)fluoranthene	15.08	252	29555m	7.577	ng	
89) Benzo(a)pyrene	15.42	252	53439	14.847	ng	98
90) Dibenzo(a,h)anthracene	16.96	278	7945m	2.504	ng	
91) Benzo(g,h,i)perylene	17.42	276	28489	9.527	ng	# 87

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

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