

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
 Data File : BF100757.D
 Acq On : 21 Nov 2017 3:28
 Operator : SJ/JU
 Sample : I6311-05
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-111717

Manual Integrations
 APPROVED

SOHIL
 11/21/2017 2:16:42 PM

Quant Time: Nov 21 06:35:21 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 15:22:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	82851	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	302827	20.00	ng	-0.02
38) Acenaphthene-d10	9.91	164	113259	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	172176	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	186816	20.00	ng	-0.01
86) Perylene-d12	15.51	264	145564	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	482328	93.32	ng	0.00
7) Phenol-d6	6.50	99	566167	88.83	ng	-0.01
23) Nitrobenzene-d5	7.43	82	337288	73.64	ng	-0.02
41) 2,4,6-Tribromophenol	10.70	330	100990	87.50	ng	-0.01
44) 2-Fluorobiphenyl	9.23	172	622116	83.64	ng	-0.01
78) Terphenyl-d14	12.98	244	456194	56.11	ng	0.00
Target Compounds						
10) Phenol	6.51	94	21669	3.239	ng	Qvalue 96
49) Dimethylphthalate	9.62	163	37477	4.334	ng	99
57) Fluorene	10.46	166	19278	2.621	ng	98
70) Phenanthrene	11.42	178	126173	13.918	ng	98
71) Anthracene	11.47	178	35698	3.881	ng	98
74) Fluoranthene	12.60	202	141768	14.756	ng	97
77) Pyrene	12.84	202	137386	10.317	ng	99
80) Benzo(a)anthracene	14.02	228	69607	6.187	ng	98
82) Chrysene	14.06	228	65603	6.022	ng	94
85) Indeno(1,2,3-cd)pyrene	16.98	276	19182	2.177	ng	93
87) Benzo(b)fluoranthene	15.07	252	57116m	6.623	ng	
88) Benzo(k)fluoranthene	15.10	252	20507m	2.517	ng	
89) Benzo(a)pyrene	15.44	252	40751	5.330	ng	94
91) Benzo(a,h,i)perylene	17.43	276	18840	3.078	ng	# 90

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

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