

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
 Data File : BF100143.D
 Acq On : 3 Nov 2017 19:10
 Operator : SJ/JU
 Sample : I6093-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B8-B9-COMP-4-12

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:18:56 PM

Quant Time: Nov 03 23:23:34 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	103146	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	427920	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	195152	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	327625	20.00	ng	0.00
75) Chrysene-d12	13.95	240	180050	20.00	ng	0.00
86) Perylene-d12	15.41	264	168575	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	643760	97.99	ng	0.00
7) Phenol-d6	6.42	99	761414	97.74	ng	0.00
23) Nitrobenzene-d5	7.34	82	436249	65.63	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	164436	92.46	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	769203	60.81	ng	0.00
78) Terphenyl-d14	12.88	244	590677	71.69	ng	0.00
Target Compounds						
10) Phenol	6.43	94	47231	5.522	ng	93
49) Dimethylphthalate	9.52	163	154314	9.985	ng	100
70) Phenanthrene	11.32	178	121663	6.580	ng	98
74) Fluoranthene	12.52	202	176251	9.173	ng	98
77) Pyrene	12.74	202	136706	9.985	ng	98
80) Benzo(a)anthracene	13.94	228	39167	3.432	ng	96
82) Chrysene	13.98	228	47022	4.382	ng	96
85) Indeno(1,2,3-cd)pyrene	16.90	276	25803	2.619	ng	96
87) Benzo(b)fluoranthene	14.99	252	52005m	4.886	ng	
89) Benzo(a)pyrene	15.36	252	33138	3.466	ng	# 94
91) Benzo(g,h,i)perylene	17.35	276	24566	2.955	ng	# 90

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

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