

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100095.D
 Acq On : 2 Nov 2017 19:55
 Operator : SJ/JU
 Sample : I6233-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3(6-8)

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:19:45 AM

Quant Time: Nov 02 20:31:47 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	88991	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	357900	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	156125	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	252330	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	149109	20.00	ng	-0.02
86) Perylene-d12	15.41	264	152193	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	584111	103.05	ng	0.00
7) Phenol-d6	6.42	99	707887	105.32	ng	-0.01
23) Nitrobenzene-d5	7.34	82	397888	71.57	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	148285	104.22	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	786678	77.74	ng	0.00
78) Terphenyl-d14	12.88	244	537851	78.83	ng	-0.02
Target Compounds						
48) Acenaphthylene	9.67	152	49260	3.118	ng	Qvalue 100
49) Dimethylphthalate	9.52	163	135977	10.998	ng	99
70) Phenanthrene	11.32	178	252296	17.717	ng	98
71) Anthracene	11.37	178	97945	6.776	ng	99
74) Fluoranthene	12.52	202	455054	30.749	ng	99
77) Pyrene	12.74	202	370679	32.693	ng	99
80) Benzo(a)anthracene	13.94	228	192645	20.380	ng	99
82) Chrysene	13.97	228	148464	16.707	ng	98
85) Indeno(1,2,3-cd)pyrene	16.88	276	83780	10.270	ng	92
87) Benzo(b)fluoranthene	14.99	252	190620m	19.835	ng	
88) Benzo(k)fluoranthene	15.01	252	69084m	7.623	ng	
89) Benzo(a)pyrene	15.35	252	145761	16.885	ng	98
90) Dibenzo(a,h)anthracene	16.87	278	20168	2.629	ng	# 78
91) Benzo(a,h,i)perylene	17.32	276	68366	9.110	ng	92

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

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