

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF120117\
 Data File : BF101065.D
 Acq On : 1 Dec 2017 16:11
 Operator : SJ/JU
 Sample : I6300-07 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-112917

Manual Integrations
 APPROVED

Sohil
 12/4/2017 11:56:41 AM

Quant Time: Dec 04 00:33:17 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 15:52:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	51207	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	196544	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	91534	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	162094	20.00	ng	0.00
75) Chrysene-d12	14.02	240	108806	20.00	ng	0.00
86) Perylene-d12	15.49	264	119951	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	63054	20.35	ng	0.00
7) Phenol-d6	6.47	99	76626	20.37	ng	-0.01
23) Nitrobenzene-d5	7.40	82	44786	13.59	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	13980	12.71	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	104409	15.61	ng	-0.01
78) Terphenyl-d14	12.95	244	72745	14.62	ng	-0.01
Target Compounds						Qvalue
51) Acenaphthene	9.92	154	16040	2.737	ng	98
57) Fluorene	10.43	166	23797	3.466	ng	97
70) Phenanthrene	11.40	178	209816	23.505	ng	98
71) Anthracene	11.45	178	47925	5.305	ng	98
74) Fluoranthene	12.59	202	243455	24.604	ng	96
77) Pyrene	12.82	202	220770	29.405	ng	99
80) Benzo(a)anthracene	14.00	228	103845	15.116	ng	98
82) Chrysene	14.04	228	90477	14.181	ng	96
85) Indeno(1,2,3-cd)pyrene	16.96	276	42842	7.553	ng	# 89
87) Benzo(b)fluoranthene	15.06	252	114191m	15.894	ng	
88) Benzo(k)fluoranthene	15.09	252	41006m	5.793	ng	
89) Benzo(a)pyrene	15.43	252	87923	13.461	ng	95
91) Benzo(a,h,i)perylene	17.41	276	40719	7.503	ng	# 91

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

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