

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\
 Data File : BF101181.D
 Acq On : 6 Dec 2017 21:19
 Operator : SJ/JU
 Sample : I6640-01 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11(0-2)-20171129

Manual Integrations
 APPROVED

Sohil
 12/7/2017 11:36:06 AM

Quant Time: Dec 07 01:08:28 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 04 13:07:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	43479	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	162010	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	67553	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	110903	20.00	ng	0.00
75) Chrysene-d12	14.02	240	102199	20.00	ng	0.00
86) Perylene-d12	15.50	264	95651	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	136235	51.78	ng	0.00
7) Phenol-d6	6.47	99	156871	49.11	ng	0.00
23) Nitrobenzene-d5	7.41	82	94100	34.65	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	31787	39.17	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	204704	41.46	ng	0.00
78) Terphenyl-d14	12.96	244	147030	31.47	ng	0.00
Target Compounds						
70) Phenanthrene	11.40	178	85452	13.992	ng	98
71) Anthracene	11.45	178	17911	2.898	ng	99
74) Fluoranthene	12.59	202	142396	21.034	ng	94
77) Pyrene	12.82	202	136711	19.386	ng	97
80) Benzo(a)anthracene	14.01	228	77025	11.937	ng	98
82) Chrysene	14.04	228	68268	11.392	ng	98
85) Indeno(1,2,3-cd)pyrene	16.99	276	30825	5.786	ng	90
87) Benzo(b)fluoranthene	15.07	252	88081m	15.374	ng	
88) Benzo(k)fluoranthene	15.09	252	25859m	4.581	ng	
89) Benzo(a)pyrene	15.44	252	58147	11.164	ng	96
91) Benzo(g,h,i)perylene	17.44	276	30854	7.130	ng	# 91

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

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