

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
 Data File : BF107892.D
 Acq On : 1 Aug 2018 16:40
 Operator : JU/SJ
 Sample : J3826-05 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-073118

Manual Integrations
 APPROVED

Sohil
 8/2/2018 9:17:12 AM

Quant Time: Aug 01 17:50:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	94851	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	430138	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	175151	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	357937	20.00	ng	0.00
75) Chrysene-d12	14.15	240	393286	20.00	ng	0.00
86) Perylene-d12	15.69	264	373861	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	97285m	17.43	ng	0.02
7) Phenol-d6	6.62	99	142313m	19.35	ng	0.02
23) Nitrobenzene-d5	7.57	82	83320	11.16	ng	0.03
41) 2,4,6-Tribromophenol	10.80	330	38157	16.03	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	183571	14.60	ng	0.00
78) Terphenyl-d14	13.08	244	238067	13.13	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	26908	2.046	ng	98
70) Phenanthrene	11.52	178	196480	11.560	ng	98
71) Anthracene	11.58	178	76177	3.886	ng	94
74) Fluoranthene	12.72	202	353064	17.682	ng	96
77) Pyrene	12.95	202	309632	10.933	ng	97
80) Benzo(a)anthracene	14.14	228	162164	7.279	ng	97
82) Chrysene	14.18	228	136379	5.676	ng	97
85) Indeno(1,2,3-cd)pyrene	17.28	276	44877	2.455	ng	# 89
87) Benzo(b)fluoranthene	15.23	252	133057m	7.216	ng	
88) Benzo(k)fluoranthene	15.25	252	80953m	3.576	ng	
89) Benzo(a)pyrene	15.62	252	104297	5.479	ng	93
91) Benzo(a,h,i)perylene	17.76	276	40730	2.535	ng	# 78

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

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