

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
 Data File : BF112969.D
 Acq On : 11 Mar 2019 21:55
 Operator : JU/SJ
 Sample : K1691-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-030819-A

Manual Integrations
 APPROVED

Sohil
 3/12/2019 4:57:31 PM

Quant Time: Mar 12 16:08:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	186981	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	790779	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	439577	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	905894	20.00	ng	0.00
76) Chrysene-d12	13.86	240	623092	20.00	ng	-0.01
87) Perylene-d12	15.27	264	590371	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	505645	42.07	ng	-0.01
7) Phenol-d6	6.36	99	676435	44.80	ng	0.00
23) Nitrobenzene-d5	7.29	82	387743	29.34	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	261521	56.04	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	789395	24.20	ng	-0.01
79) Terphenyl-d14	12.82	244	847548	26.37	ng	-0.01
Target Compounds						
58) Fluorene	10.30	166	61279	2.137	ng	99
71) Phenanthrene	11.26	178	419198	9.301	ng	99
72) Anthracene	11.32	178	140556	2.979	ng	97
75) Fluoranthene	12.44	202	625066	12.944	ng	96
78) Pyrene	12.67	202	517254	9.847	ng	98
81) Benzo(a)anthracene	13.85	228	225306	5.217	ng	99
83) Chrysene	13.89	228	196485	4.645	ng	96
86) Indeno(1,2,3-cd)pyrene	16.62	276	86558m	2.400	ng	
88) Benzo(b)fluoranthene	14.87	252	236829m	6.202	ng	
89) Benzo(k)fluoranthene	14.90	252	101243m	2.927	ng	
90) Benzo(a)pyrene	15.21	252	179491	5.314	ng	98
92) Benzo(a,h,i)perylene	17.03	276	75634	2.613	ng	94

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

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