

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080318\
 Data File : BF107970.D
 Acq On : 4 Aug 2018 6:21
 Operator : JU/SJ
 Sample : J4262-01DL 2X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-04DL

Manual Integrations
 APPROVED

Sohil
 8/6/2018 5:06:14 PM

Quant Time: Aug 06 04:10:36 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	6.96	152	74416	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	348283	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	143497	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	257048	20.00	ng	0.00
75) Chrysene-d12	14.15	240	255232	20.00	ng	0.00
86) Perylene-d12	15.68	264	193182	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	70047	16.00	ng	0.02
7) Phenol-d6	6.64	99	86968	15.07	ng	0.03
23) Nitrobenzene-d5	7.52	82	232667	38.49	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	103875	53.27	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	518275	50.33	ng	-0.01
78) Terphenyl-d14	13.08	244	521328	44.29	ng	0.00
Target Compounds						
31) Naphthalene	8.26	128	854239	52.764	ng	100
37) 2-Methylnaphthalene	8.95	142	75397	7.414	ng	95
48) Acenaphthylene	9.86	152	39986	2.619	ng	# 93
51) Acenaphthene	10.02	154	135670	15.298	ng	100
54) Dibenzofuran	10.21	168	85671	6.250	ng	96
57) Fluorene	10.55	166	89902	8.439	ng	97
70) Phenanthrene	11.52	178	104643	8.573	ng	99
71) Anthracene	11.58	178	35398m	2.514	ng	
72) Carbazole	11.74	167	181363	13.036	ng	98

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

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