

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
 Data File : BF107925.D
 Acq On : 2 Aug 2018 22:28
 Operator : JU/SJ
 Sample : J4285-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KG-PIT-3

Manual Integrations
 APPROVED

Sohil
 8/3/2018 12:53:35 PM

Quant Time: Aug 03 04:39:52 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	6.95	152	117496	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	460485	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	166440	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	290840	20.00	ng	-0.01
75) Chrysene-d12	14.16	240	260966	20.00	ng	0.00
86) Perylene-d12	15.69	264	229333	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	597164	86.38	ng	0.00
7) Phenol-d6	6.60	99	785761	86.24	ng	0.00
23) Nitrobenzene-d5	7.52	82	437214	54.70	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	162278	71.75	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	780361	65.33	ng	-0.01
78) Terphenyl-d14	13.08	244	586671	48.75	ng	0.00
Target Compounds						
31) Naphthalene	8.26	128	52074	2.433	ng	99
49) Dimethylphthalate	9.70	163	113125	9.050	ng	100
70) Phenanthrene	11.52	178	58858	4.262	ng	# 85
74) Fluoranthene	12.72	202	92382	5.694	ng	75
77) Pyrene	12.95	202	95195	5.066	ng	# 77
80) Benzo(a)anthracene	14.15	228	43685	2.955	ng	# 53
82) Chrysene	14.19	228	64700	4.058	ng	# 80
85) Indeno(1,2,3-cd)pyrene	17.27	276	36718	3.028	ng	# 83
87) Benzo(b)fluoranthene	15.24	252	68189m	6.029	ng	
88) Benzo(k)fluoranthene	15.25	252	47051m	3.388	ng	
89) Benzo(a)pyrene	15.62	252	37561	3.217	ng	# 52
91) Benzo(a,h,i)perylene	17.75	276	34764	3.527	ng	# 88

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

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