

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108096.D
 Acq On : 10 Aug 2018 20:51
 Operator : JU/SJ
 Sample : J4272-15 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-080918-H

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:52:20 PM

Quant Time: Aug 11 02:06:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	285882	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1026795	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	429049	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	708702	20.00	ng	0.00
75) Chrysene-d12	14.22	240	583414	20.00	ng	0.00
86) Perylene-d12	15.79	264	391894	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	784435	49.68	ng	0.00
7) Phenol-d6	6.63	99	1045690	49.83	ng	0.00
23) Nitrobenzene-d5	7.58	82	647999	35.22	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	192660	45.02	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1139443	42.64	ng	0.00
78) Terphenyl-d14	13.15	244	989295	43.11	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.77	163	107584	3.600	ng	Qvalue # 99
70) Phenanthrene	11.59	178	193477	5.788	ng	98
71) Anthracene	11.64	178	71206	2.078	ng	98
74) Fluoranthene	12.79	202	450297	12.343	ng	93
77) Pyrene	13.02	202	421148	11.402	ng	100
80) Benzo(a)anthracene	14.21	228	212990	6.361	ng	96
82) Chrysene	14.25	228	158935	5.178	ng	97
85) Indeno(1,2,3-cd)pyrene	17.41	276	59630	2.242	ng	# 90
87) Benzo(b)fluoranthene	15.32	252	168527m	7.197	ng	
88) Benzo(k)fluoranthene	15.35	252	63016m	2.927	ng	
89) Benzo(a)pyrene	15.72	252	113507	5.413	ng	# 92
91) Benzo(g,h,i)perylene	17.91	276	58386	3.417	ng	# 93

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

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