

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108095.D
 Acq On : 10 Aug 2018 20:23
 Operator : JU/SJ
 Sample : J4272-13 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:52:18 PM

Quant Time: Aug 11 01:45:13 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	293910	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	1063231	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	468870	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	763997	20.00	ng	0.00
75) Chrysene-d12	14.22	240	630561	20.00	ng	0.00
86) Perylene-d12	15.79	264	436482	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	804902	49.58	ng	0.00
7) Phenol-d6	6.63	99	1080696	50.10	ng	0.00
23) Nitrobenzene-d5	7.58	82	673123	35.33	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	212033	45.34	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1211781	41.49	ng	0.00
78) Terphenyl-d14	13.15	244	1020901	41.17	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.77	163	96458	2.953	ng	99
74) Fluoranthene	12.78	202	146189	3.717	ng	96
77) Pyrene	13.02	202	139918	3.505	ng	99
80) Benzo(a)anthracene	14.21	228	78940	2.181	ng	93
82) Chrysene	14.25	228	73783	2.224	ng	98
87) Benzo(b)fluoranthene	15.32	252	80795m	3.098	ng	
88) Benzo(k)fluoranthene	15.32	252	98863	4.124	ng	# 84
89) Benzo(a)pyrene	15.72	252	50408	2.158	ng	# 78

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

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