

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108094.D
 Acq On : 10 Aug 2018 19:55
 Operator : JU/SJ
 Sample : J4272-11 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 8/13/2018 3:52:15 PM

Quant Time: Aug 11 01:30:14 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	268116	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	944124	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	413287	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	697696	20.00	ng	0.00
75) Chrysene-d12	14.22	240	614361	20.00	ng	0.00
86) Perylene-d12	15.79	264	412167	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	741019	50.04	ng	0.00
7) Phenol-d6	6.63	99	996921	50.66	ng	0.00
23) Nitrobenzene-d5	7.58	82	616583	36.44	ng	0.00
41) 2,4,6-Tribromophenol	10.86	330	196113	47.57	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1031552	40.07	ng	0.00
78) Terphenyl-d14	13.15	244	947305	39.20	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.77	163	81949	2.847	ng	99
70) Phenanthrene	11.59	178	104950	3.189	ng	100
74) Fluoranthene	12.78	202	235255	6.550	ng	97
77) Pyrene	13.02	202	221235	5.688	ng	99
80) Benzo(a)anthracene	14.21	228	119752	3.396	ng	96
82) Chrysene	14.25	228	101940	3.154	ng	96
87) Benzo(b)fluoranthene	15.31	252	107732m	4.374	ng	
89) Benzo(a)pyrene	15.72	252	68953	3.127	ng	# 89

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

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