

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081318\
 Data File : BF108113.D
 Acq On : 13 Aug 2018 13:55
 Operator : JU/SJ
 Sample : J4272-11 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 8/14/2018 2:36:24 PM

Quant Time: Aug 13 16:11:52 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	264471	20.00	ng	0.00
21) Naphthalene-d8	8.31	136	993575	20.00	ng	0.00
38) Acenaphthene-d10	10.07	164	487270	20.00	ng	0.00
63) Phenanthrene-d10	11.57	188	857961	20.00	ng	0.01
75) Chrysene-d12	14.23	240	667215	20.00	ng	0.01
86) Perylene-d12	15.82	264	693969	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	737678	50.50	ng	0.00
7) Phenol-d6	6.63	99	1023715	52.74	ng	0.00
23) Nitrobenzene-d5	7.58	82	644995	36.22	ng	0.00
41) 2,4,6-Tribromophenol	10.87	330	245443	50.50	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1102697	36.33	ng	0.00
78) Terphenyl-d14	13.16	244	972041	37.04	ng	0.01
Target Compounds						
49) Dimethylphthalate	9.77	163	91491	2.695	ng	Qvalue # 99
70) Phenanthrene	11.59	178	127523	3.151	ng	97
74) Fluoranthene	12.79	202	250923	5.682	ng	97
77) Pyrene	13.02	202	236536	5.600	ng	99
80) Benzo(a)anthracene	14.22	228	129710	3.387	ng	97
82) Chrysene	14.25	228	108793	3.099	ng	97
87) Benzo(b)fluoranthene	15.34	252	148674m	3.585	ng	
89) Benzo(a)pyrene	15.74	252	114179	3.075	ng	92

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

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