

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106356.D
 Acq On : 12 Jun 2018 20:30
 Operator : JU/SJ
 Sample : J3354-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CAS1-SB-102-103-1-27

Manual Integrations
 APPROVED

Sohil
 6/13/2018 4:56:14 PM

Quant Time: Jun 13 03:26:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	146879	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	495243	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	199676	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	392401	20.00	ng	0.00
75) Chrysene-d12	14.15	240	427038	20.00	ng	0.00
86) Perylene-d12	15.69	264	311798	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1058774	122.00	ng	0.02
7) Phenol-d6	6.60	99	1271841	116.00	ng	0.01
23) Nitrobenzene-d5	7.54	82	808595	96.05	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	303092	157.76	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1243083	129.58	ng	0.00
78) Terphenyl-d14	13.09	244	1508227	87.72	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.72	163	138382	9.700	ng	99
70) Phenanthrene	11.53	178	70503	3.669	ng	98
74) Fluoranthene	12.72	202	179237	8.746	ng	96
77) Pyrene	12.95	202	234637	8.776	ng	99
80) Benzo(a)anthracene	14.14	228	104762	4.122	ng	99
82) Chrysene	14.18	228	97836m	4.217	ng	
87) Benzo(b)fluoranthene	15.23	252	77004m	4.086	ng	
89) Benzo(a)pyrene	15.62	252	59502	3.510	ng	95

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

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