

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106688.D
 Acq On : 22 Jun 2018 3:34
 Operator : JU/SJ
 Sample : J3603-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26KV-TP-1

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:46 AM

Quant Time: Jun 22 04:44:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	56709	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	193099	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	87744	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	178759	20.00	ng	0.00
75) Chrysene-d12	14.15	240	148701	20.00	ng	0.00
86) Perylene-d12	15.68	264	111522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	285176	85.15	ng	0.00
7) Phenol-d6	6.60	99	344690	82.42	ng	0.00
23) Nitrobenzene-d5	7.52	82	196779	56.63	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	82780	81.40	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	282661	48.13	ng	0.00
78) Terphenyl-d14	13.08	244	330090	53.77	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	31662	4.811	ng	97
70) Phenanthrene	11.52	178	35638	4.133	ng	99
74) Fluoranthene	12.71	202	59097	6.219	ng	98
77) Pyrene	12.94	202	60823	6.203	ng	99
80) Benzo(a)anthracene	14.14	228	32618	3.599	ng	# 86
82) Chrysene	14.17	228	29150	3.496	ng	# 95
87) Benzo(b)fluoranthene	15.23	252	26002m	3.753	ng	
89) Benzo(a)pyrene	15.62	252	20181	3.277	ng	# 89
91) Benzo(g,h,i)perylene	17.72	276	12624m	2.475	ng	

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

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