

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106702.D
 Acq On : 22 Jun 2018 12:10
 Operator : JU/SJ
 Sample : J3603-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26KV-TP-8

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:16:36 PM

Quant Time: Jun 22 12:40:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 22 11:34:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	64940	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	253062	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	133738	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	262355	20.00	ng	0.00
75) Chrysene-d12	14.15	240	199472	20.00	ng	0.00
86) Perylene-d12	15.69	264	207613	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	408182	106.43	ng	0.00
7) Phenol-d6	6.61	99	524225	109.46	ng	0.00
23) Nitrobenzene-d5	7.52	82	352434	77.40	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	168537	108.73	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	510548	60.34	ng	0.00
78) Terphenyl-d14	13.08	244	533338	64.77	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	34681	3.458	ng	99
70) Phenanthrene	11.52	178	35042	2.769	ng	98
74) Fluoranthene	12.71	202	43847	3.144	ng	98
77) Pyrene	12.94	202	39914	3.034	ng	98
80) Benzo(a)anthracene	14.14	228	25063	2.062	ng	84
82) Chrysene	14.17	228	31941	2.856	ng	93
87) Benzo(b)fluoranthene	15.23	252	30479m	2.363	ng	
89) Benzo(a)pyrene	15.62	252	23633	2.061	ng	98

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

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