

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106701.D
 Acq On : 22 Jun 2018 11:44
 Operator : JU/SJ
 Sample : J3603-15
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26KV-TP-9

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:16:33 PM

Quant Time: Jun 22 12:31:30 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	67281	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	258136	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	136075	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	268374	20.00	ng	0.00
75) Chrysene-d12	14.15	240	195129	20.00	ng	0.00
86) Perylene-d12	15.69	264	210586	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	393016	98.91	ng	0.00
7) Phenol-d6	6.61	99	516296	104.05	ng	0.00
23) Nitrobenzene-d5	7.52	82	379299	81.66	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	135890	86.16	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	532612	62.31	ng	0.00
78) Terphenyl-d14	13.08	244	481636	59.79	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	82555	8.089	ng	98
70) Phenanthrene	11.52	178	100178	7.739	ng	98
74) Fluoranthene	12.71	202	160630	11.259	ng	96
77) Pyrene	12.94	202	153965	11.966	ng	99
80) Benzo(a)anthracene	14.14	228	87403	7.349	ng	99
82) Chrysene	14.17	228	81621	7.460	ng	97
85) Indeno(1,2,3-cd)pyrene	17.25	276	41955	4.519	ng	# 89
87) Benzo(b)fluoranthene	15.23	252	97264m	7.435	ng	
88) Benzo(k)fluoranthene	15.26	252	26005m	2.088	ng	
89) Benzo(a)pyrene	15.62	252	72065m	6.197	ng	
91) Benzo(g,h,i)perylene	17.73	276	42889	4.452	ng	# 91

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

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