

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062518\
 Data File : BF106786.D
 Acq On : 25 Jun 2018 21:55
 Operator : JU/SJ
 Sample : J3241-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-062218-C

Manual Integrations
 APPROVED

Sohil
 6/27/2018 10:55:41 AM

Quant Time: Jun 26 03:45:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	45381	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	164287	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	70920	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	123795	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	133256	20.00	ng	-0.02
86) Perylene-d12	15.69	264	102938	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	174131	64.97	ng	0.00
7) Phenol-d6	6.60	99	213421	63.77	ng	-0.02
23) Nitrobenzene-d5	7.52	82	128998	43.64	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	52012	63.28	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	246785	53.42	ng	-0.01
78) Terphenyl-d14	13.07	244	269068	48.91	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.71	163	40612	7.635	ng	Qvalue # 98
70) Phenanthrene	11.51	178	61957	10.377	ng	97
71) Anthracene	11.57	178	19966	3.276	ng	99
74) Fluoranthene	12.71	202	130928	19.895	ng	97
77) Pyrene	12.94	202	126315	14.375	ng	98
80) Benzo(a)anthracene	14.14	228	75063	9.242	ng	98
82) Chrysene	14.17	228	68382	9.152	ng	97
85) Indeno(1,2,3-cd)pyrene	17.26	276	27129	4.279	ng	# 89
87) Benzo(b)fluoranthene	15.23	252	80213m	12.544	ng	
88) Benzo(k)fluoranthene	15.25	252	24388m	4.005	ng	
89) Benzo(a)pyrene	15.62	252	52480	9.232	ng	93
91) Benzo(a,h,i)perylene	17.74	276	29678	6.303	ng	# 87

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

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