

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
 Data File : BF107042.D
 Acq On : 2 Jul 2018 14:13
 Operator : JU/SJ
 Sample : J3774-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6(2-4)

Manual Integrations
 APPROVED

Sohil
 7/3/2018 5:29:46 PM

Quant Time: Jul 03 08:55:26 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 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	63632	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	235018	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	112040	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	206061	20.00	ng	0.00
75) Chrysene-d12	14.15	240	183590	20.00	ng	-0.01
86) Perylene-d12	15.71	264	149337	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	359625	95.70	ng	0.00
7) Phenol-d6	6.60	99	454191	96.78	ng	-0.01
23) Nitrobenzene-d5	7.52	82	280757	66.39	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	137026	105.52	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	550512	82.78	ng	-0.01
78) Terphenyl-d14	13.08	244	576694	76.09	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	48411	5.761	ng	97
70) Phenanthrene	11.52	178	50181	5.049	ng	99
74) Fluoranthene	12.71	202	66606	6.080	ng	96
77) Pyrene	12.94	202	110867	9.158	ng	98
80) Benzo(a)anthracene	14.14	228	80918	7.232	ng	97
82) Chrysene	14.18	228	68348	6.640	ng	100
85) Indeno(1,2,3-cd)pyrene	17.29	276	45916	5.256	ng	# 86
87) Benzo(b)fluoranthene	15.25	252	93829m	10.114	ng	
88) Benzo(k)fluoranthene	15.27	252	27977m	3.167	ng	
89) Benzo(a)pyrene	15.64	252	102514m	12.431	ng	
90) Dibenzo(a,h)anthracene	17.29	278	15218	2.177	ng	94
91) Benzo(a,h,i)perylene	17.77	276	51904	7.598	ng	# 89

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

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