

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062818\
 Data File : BF106936.D
 Acq On : 28 Jun 2018 22:17
 Operator : JU/SJ
 Sample : J3242-13 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-062618-G

Manual Integrations
 APPROVED

Sohil
 6/29/2018 2:30:53 PM

Quant Time: Jun 29 02:13:17 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	6.95	152	65585	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	225955	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	90945	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	166752	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	215785	20.00	ng	-0.03
86) Perylene-d12	15.67	264	191651	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	173597	44.82	ng	-0.01
7) Phenol-d6	6.59	99	205188	42.42	ng	-0.02
23) Nitrobenzene-d5	7.51	82	121694	29.93	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	45055	42.74	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	225814	33.00	ng	-0.02
78) Terphenyl-d14	13.07	244	274706	30.84	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	38549	5.652	ng	99
70) Phenanthrene	11.51	178	19948	2.480	ng	98
74) Fluoranthene	12.70	202	46218	5.214	ng	95
77) Pyrene	12.94	202	46746	3.285	ng	97
80) Benzo(a)anthracene	14.12	228	35648	2.711	ng	97
82) Chrysene	14.16	228	31731	2.623	ng	97
87) Benzo(b)fluoranthene	15.21	252	36934m	3.102	ng	
89) Benzo(a)pyrene	15.60	252	23605	2.230	ng	# 91

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

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