

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062818\
 Data File : BF106932.D
 Acq On : 28 Jun 2018 20:29
 Operator : JU/SJ
 Sample : J3242-15 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 6/29/2018 2:30:48 PM

Quant Time: Jun 29 01:48:48 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	79219	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	304183	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	143495	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	232476	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	224059	20.00	ng	-0.03
86) Perylene-d12	15.67	264	223437	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	252739	54.02	ng	-0.01
7) Phenol-d6	6.59	99	316440	54.16	ng	-0.02
23) Nitrobenzene-d5	7.51	82	191787	35.04	ng	-0.02
41) 2,4,6-Tribromophenol	10.78	330	85311	51.29	ng	-0.02
44) 2-Fluorobiphenyl	9.31	172	396634	38.76	ng	-0.02
78) Terphenyl-d14	13.07	244	355665	38.45	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.70	163	78346	7.280	ng	98
70) Phenanthrene	11.51	178	24862	2.217	ng	98
74) Fluoranthene	12.70	202	52355	4.236	ng	94
77) Pyrene	12.94	202	50337	3.407	ng	99
80) Benzo(a)anthracene	14.12	228	30812	2.256	ng	97
82) Chrysene	14.16	228	27378	2.179	ng	96
87) Benzo(b)fluoranthene	15.21	252	41937m	3.021	ng	
89) Benzo(a)pyrene	15.60	252	28058	2.274	ng	95

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

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