

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071318\
 Data File : BF107368.D
 Acq On : 13 Jul 2018 15:42
 Operator : JU/SJ
 Sample : J3825-03 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-071118-B

Manual Integrations
 APPROVED

Sohil
 7/16/2018 3:22:23 PM

Quant Time: Jul 13 18:02:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 13 11:54:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	49924	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	174164	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	78167	20.00	ng	0.00
63) Phenanthrene-d10	11.47	188	164942	20.00	ng	0.00
75) Chrysene-d12	14.14	240	174658	20.00	ng	0.00
86) Perylene-d12	15.68	264	120271	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	136313	46.23	ng	0.00
7) Phenol-d6	6.58	99	154959	42.09	ng	0.00
23) Nitrobenzene-d5	7.50	82	91242	29.11	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	39633	43.75	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	175737	28.20	ng	0.00
78) Terphenyl-d14	13.05	244	246298	34.16	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	24257	4.138	ng	97
70) Phenanthrene	11.50	178	41088	5.165	ng	97
74) Fluoranthene	12.69	202	72017	8.213	ng	96
77) Pyrene	12.93	202	84492	7.336	ng	97
80) Benzo(a)anthracene	14.12	228	41891	3.935	ng	94
82) Chrysene	14.16	228	36574	3.735	ng	97
87) Benzo(b)fluoranthene	15.22	252	33822m	4.527	ng	
89) Benzo(a)pyrene	15.61	252	24220	3.647	ng	# 87
91) Benzo(g,h,i)perylene	17.77	276	17457	3.173	ng	# 91

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

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