

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071618\
 Data File : BF107416.D
 Acq On : 16 Jul 2018 21:34
 Operator : JU/SJ
 Sample : J3820-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-071218

Manual Integrations
 APPROVED

Sohil
 7/17/2018 4:53:05 PM

Quant Time: Jul 17 02:01:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 16 17:24:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	165022	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	604677	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	286803	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	539357	20.00	ng	0.00
75) Chrysene-d12	14.16	240	488020	20.00	ng	0.00
86) Perylene-d12	15.69	264	453142	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	911820	83.92	ng	0.00
7) Phenol-d6	6.60	99	1188046	84.57	ng	0.00
23) Nitrobenzene-d5	7.54	82	841983	74.25	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	226560	89.91	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	1333481	76.43	ng	0.00
78) Terphenyl-d14	13.10	244	1226886	62.46	ng	0.00
Target Compounds						
10) Phenol	6.61	94	38332	2.545	ng	72
49) Dimethylphthalate	9.72	163	251987	11.058	ng	98
70) Phenanthrene	11.54	178	231023	8.595	ng	98
74) Fluoranthene	12.73	202	251796	8.242	ng	92
77) Pyrene	12.96	202	355371	10.708	ng	94
80) Benzo(a)anthracene	14.15	228	141231m	4.730	ng	
82) Chrysene	14.18	228	192235	6.739	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	54844	2.701	ng	96
87) Benzo(b)fluoranthene	15.23	252	144273m	5.440	ng	
88) Benzo(k)fluoranthene	15.26	252	64346m	2.486	ng	
89) Benzo(a)pyrene	15.62	252	139170	5.717	ng	99
91) Benzo(a,h,i)perylene	17.72	276	62345	3.340	ng	95

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

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