

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071318\
 Data File : BF107369.D
 Acq On : 13 Jul 2018 16:09
 Operator : JU/SJ
 Sample : J3825-05 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 7/16/2018 3:22:25 PM

Quant Time: Jul 13 18:14:38 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	43584	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	147527	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	68899	20.00	ng	0.00
63) Phenanthrene-d10	11.47	188	157298	20.00	ng	0.00
75) Chrysene-d12	14.14	240	157949	20.00	ng	0.00
86) Perylene-d12	15.68	264	107826	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	123938	48.15	ng	0.00
7) Phenol-d6	6.58	99	141097	43.90	ng	0.00
23) Nitrobenzene-d5	7.50	82	80362	30.27	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	39141	49.01	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	158245	29.19	ng	0.00
78) Terphenyl-d14	13.06	244	247378	37.94	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	20102	3.890	ng	Qvalue # 96
70) Phenanthrene	11.50	178	22033	2.904	ng	97
74) Fluoranthene	12.70	202	56981	6.814	ng	95
77) Pyrene	12.93	202	58903	5.655	ng	99
80) Benzo(a)anthracene	14.12	228	31271	3.248	ng	94
82) Chrysene	14.16	228	27256	3.078	ng	95
87) Benzo(b)fluoranthene	15.22	252	27050m	4.038	ng	
89) Benzo(a)pyrene	15.61	252	18765	3.151	ng	# 91
91) Benzo(g,h,i)perylene	17.77	276	11410	2.313	ng	# 85

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

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