

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071317\
 Data File : BF096741.D
 Acq On : 13 Jul 2017 21:38
 Operator : SJ/JU
 Sample : I4147-01 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 F-7-11RE

Manual Integrations
 APPROVED

mohammad
 7/17/2017 8:11:46 AM

Quant Time: Jul 14 03:21:08 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 14:49:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	115853	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	464331	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	168843	20.00	ng	0.00
63) Phenanthrene-d10	11.14	188	264636	20.00	ng	0.00
75) Chrysene-d12	13.77	240	195204	20.00	ng	0.00
86) Perylene-d12	15.16	264	179453	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	27430	3.56	ng	0.00
7) Phenol-d6	6.26	99	146621	15.86	ng	-0.02
23) Nitrobenzene-d5	7.19	82	106081	14.15	ng	-0.01
41) 2,4,6-Tribromophenol	10.45	330	405	0.28	ng	-0.01
44) 2-Fluorobiphenyl	8.99	172	187076	17.51	ng	0.00
78) Terphenyl-d14	12.72	244	105085	9.34	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.39	163	44719	3.517	ng	99
70) Phenanthrene	11.16	178	71971	5.002	ng	97
74) Fluoranthene	12.34	202	68061	4.942	ng	99
77) Pyrene	12.57	202	56070	3.165	ng	98
80) Benzo(a)anthracene	13.76	228	28309	2.328	ng	98
82) Chrysene	13.79	228	28952m	2.418	ng	
87) Benzo(b)fluoranthene	14.77	252	29556m	2.731	ng	

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

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