

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062117\
 Data File : BF096114.D
 Acq On : 22 Jun 2017 9:05
 Operator : SJ/MA
 Sample : I3792-01DL 25X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-061917DL

Manual Integrations
 APPROVED

mohammad
 6/22/2017 6:03:09 PM

Quant Time: Jun 22 11:10:56 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 16:05:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.30	152	68589	20.00	ng	0.00
21) Naphthalene-d8	9.33	136	307743	20.00	ng	0.00
38) Acenaphthene-d10	12.14	164	105866	20.00	ng	0.00
63) Phenanthrene-d10	14.54	188	189566	20.00	ng	0.00
75) Chrysene-d12	18.27	240	162141	20.00	ng	0.00
86) Perylene-d12	19.94	264	97290	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1765	0.41	ng	0.26
7) Phenol-d6	7.13	99	17697m	3.43	ng	0.25
23) Nitrobenzene-d5	8.36	82	6417m	1.29	ng	0.14
41) 2,4,6-Tribromophenol	13.50	330	1827	1.95	ng	0.05
44) 2-Fluorobiphenyl	11.12	172	23621	3.10	ng	0.00
78) Terphenyl-d14	16.92	244	20600	3.15	ng	0.00
Target Compounds						
70) Phenanthrene	14.57	178	183726	16.80	ng	98
71) Anthracene	14.67	178	65615	5.75	ng	99
74) Fluoranthene	16.36	202	399394	36.89	ng	98
77) Pyrene	16.67	202	339732	28.08	ng	96
80) Benzo(a)anthracene	18.26	228	143953	15.27	ng	97
82) Chrysene	18.30	228	112550	11.95	ng	98
85) Indeno(1,2,3-cd)pyrene	21.11	276	40249	4.99	ng	# 91
87) Benzo(b)fluoranthene	19.54	252	103493m	16.99	ng	
88) Benzo(k)fluoranthene	19.57	252	42589m	7.31	ng	
89) Benzo(a)pyrene	19.89	252	78847	14.08	ng	97
90) Dibenzo(a,h)anthracene	21.10	278	10401m	2.19	ng	
91) Benzo(a,h,i)perylene	21.44	276	36064	7.45	ng	100

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

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