

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051717\
 Data File : BF095201.D
 Acq On : 17 May 2017 23:52
 Operator : SJ/MA
 Sample : I3172-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-04

Manual Integrations
 APPROVED

mohammad
 5/18/2017 2:29:42 PM

Quant Time: May 18 09:14:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 04:48:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	78766	20.00	ng	0.00
21) Naphthalene-d8	9.79	136	360571	20.00	ng	0.00
38) Acenaphthene-d10	12.62	164	163817	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	272067	20.00	ng	0.00
75) Chrysene-d12	18.69	240	183919	20.00	ng	0.00
86) Perylene-d12	20.37	264	175045	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	530867	112.37	ng	0.00
7) Phenol-d6	7.31	99	706348	124.61	ng	0.00
23) Nitrobenzene-d5	8.67	82	410919	71.86	ng	0.00
41) 2,4,6-Tribromophenol	13.92	330	137539	102.52	ng	0.00
44) 2-Fluorobiphenyl	11.56	172	805165	70.74	ng	0.00
78) Terphenyl-d14	17.34	244	503182	60.26	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	12.26	163	108014	8.03	ng	99
74) Fluoranthene	16.81	202	60027	3.74	ng	97
77) Pyrene	17.11	202	58714	4.27	ng	96
80) Benzo(a)anthracene	18.68	228	55717	5.21	ng	98
82) Chrysene	18.73	228	71641	6.92	ng	96
85) Indeno(1,2,3-cd)pyrene	21.71	276	71220	8.47	ng	92
87) Benzo(b)fluoranthene	19.97	252	138445m	12.80	ng	
88) Benzo(k)fluoranthene	19.99	252	43104m	4.13	ng	
89) Benzo(a)pyrene	20.31	252	90469	9.06	ng	93
91) Benzo(g,h,i)perylene	22.09	276	64634	7.99	ng	93

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

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