

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051717\
 Data File : BF095204.D
 Acq On : 18 May 2017 1:29
 Operator : SJ/MA
 Sample : I3182-07 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-137-339-120

Manual Integrations
 APPROVED

mohammad
 5/18/2017 2:29:51 PM

Quant Time: May 18 07:01:44 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	82296	20.00	ng	0.00
21) Naphthalene-d8	9.79	136	324535	20.00	ng	0.00
38) Acenaphthene-d10	12.62	164	156940	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	208998	20.00	ng	0.00
75) Chrysene-d12	18.69	240	192262	20.00	ng	0.00
86) Perylene-d12	20.37	264	167649	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	245615	49.76	ng	0.00
7) Phenol-d6	7.32	99	319782	53.99	ng	0.00
23) Nitrobenzene-d5	8.67	82	157590	30.62	ng	0.00
41) 2,4,6-Tribromophenol	13.92	330	50146	39.02	ng	0.00
44) 2-Fluorobiphenyl	11.56	172	388366	35.62	ng	0.00
78) Terphenyl-d14	17.34	244	235545	26.98	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.26	163	62098	4.82	ng	99
70) Phenanthrene	15.06	178	55604	4.63	ng	97
74) Fluoranthene	16.81	202	101706	8.26	ng	98
77) Pyrene	17.11	202	85087	5.92	ng	95
80) Benzo(a)anthracene	18.68	228	41837	3.74	ng	96
82) Chrysene	18.72	228	42374	3.92	ng	96
87) Benzo(b)fluoranthene	19.97	252	48807m	4.71	ng	
88) Benzo(k)fluoranthene	19.99	252	23384m	2.34	ng	
89) Benzo(a)pyrene	20.31	252	37053	3.88	ng	# 89

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

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