

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\
 Data File : BF097448.D
 Acq On : 8 Aug 2017 00:09
 Operator : SJ/JU
 Sample : I4582-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S1

Manual Integrations
 APPROVED

Sohil
 8/9/2017 9:21:33 PM

Quant Time: Aug 08 06:50:14 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.41	152	99373	20.00	ng	0.00
21) Naphthalene-d8	7.69	136	353051	20.00	ng	0.00
38) Acenaphthene-d10	9.43	164	136273	20.00	ng	0.00
63) Phenanthrene-d10	10.91	188	200492	20.00	ng	0.00
75) Chrysene-d12	13.53	240	202910	20.00	ng	0.00
86) Perylene-d12	14.87	264	141886	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.00	112	405451	61.51	ng	0.00
7) Phenol-d6	6.08	99	471762	62.69	ng	0.00
23) Nitrobenzene-d5	6.98	82	285431	47.43	ng	-0.01
41) 2,4,6-Tribromophenol	10.23	330	61649	57.26	ng	0.00
44) 2-Fluorobiphenyl	8.77	172	427170	53.20	ng	0.00
78) Terphenyl-d14	12.49	244	311605	31.31	ng	0.00
Target Compounds						
10) Phenol	6.10	94	19490	2.183	ng	# 77
49) Dimethylphthalate	9.17	163	90364	8.732	ng	# 99
70) Phenanthrene	10.93	178	29470	2.743	ng	99
74) Fluoranthene	12.12	202	70168	6.668	ng	99
77) Pyrene	12.34	202	62567	3.766	ng	99
80) Benzo(a)anthracene	13.52	228	39298	2.993	ng	97
82) Chrysene	13.56	228	33700	2.629	ng	95
87) Benzo(b)fluoranthene	14.53	252	41459m	4.861	ng	
89) Benzo(a)pyrene	14.83	252	25174	3.225	ng	# 93

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

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