

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072817\
 Data File : BF097149.D
 Acq On : 28 Jul 2017 12:47
 Operator : SJ/JU
 Sample : I4397-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-1A-4.0-4.5-072117

Manual Integrations
 APPROVED

Sohil
 7/31/2017 6:51:36 PM

Quant Time: Jul 28 13:56:18 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	115380	20.00	ng	-0.01
21) Naphthalene-d8	7.77	136	502072	20.00	ng	-0.02
38) Acenaphthene-d10	9.52	164	199608	20.00	ng	-0.02
63) Phenanthrene-d10	10.99	188	285758	20.00	ng	-0.02
75) Chrysene-d12	13.61	240	214134	20.00	ng	-0.02
86) Perylene-d12	14.96	264	190435	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.09	112	791646	103.43	ng	0.00
7) Phenol-d6	6.15	99	921624	105.48	ng	-0.02
23) Nitrobenzene-d5	7.06	82	577318	67.46	ng	-0.02
41) 2,4,6-Tribromophenol	10.30	330	156204	99.05	ng	-0.02
44) 2-Fluorobiphenyl	8.85	172	927391	78.85	ng	-0.02
78) Terphenyl-d14	12.57	244	558889	53.21	ng	-0.01
Target Compounds						
10) Phenol	6.16	94	39883	3.848	ng	97
49) Dimethylphthalate	9.25	163	203661	13.436	ng	97
74) Fluoranthene	12.19	202	38533	2.569	ng	99
77) Pyrene	12.42	202	36962	2.108	ng	95
87) Benzo(b)fluoranthene	14.60	252	26811m	2.342	ng	

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

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