

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041917\
 Data File : BF094532.D
 Acq On : 19 Apr 2017 22:49
 Operator : SJ/MA
 Sample : I2749-03 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-V6-NSW-02

Quant Time: Apr 20 08:08:10 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.77	152	131097	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	535821	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	214197	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	404812	20.00	ng	0.00
75) Chrysene-d12	14.46	240	293970	20.00	ng	0.00
86) Perylene-d12	16.08	264	282003	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.31	112	177619	22.48	ng	0.00
7) Phenol-d6	5.38	99	224127	24.06	ng	0.00
23) Nitrobenzene-d5	6.41	82	129661	14.94	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	42017	25.08	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	261510	17.96	ng	0.00
78) Terphenyl-d14	13.19	244	194406	17.68	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.08	163	70399	4.41	ng	99
70) Phenanthrene	11.24	178	78115	3.43	ng	99
74) Fluoranthene	12.70	202	97690	4.13	ng	98
77) Pyrene	12.98	202	104708	5.10	ng	96
80) Benzo(a)anthracene	14.45	228	38980	2.28	ng	96
82) Chrysene	14.49	228	39508	2.53	ng	94
87) Benzo(b)fluoranthene	15.68	252	44691m	2.59	ng	
89) Benzo(a)pyrene	16.02	252	36397	2.27	ng	# 97

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

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