

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041917\
 Data File : BF094527.D
 Acq On : 19 Apr 2017 20:31
 Operator : SJ/MA
 Sample : I2749-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-V6-NSW-01

Quant Time: Apr 20 04:50:46 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	132827	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	505093	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	205167	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	397844	20.00	ng	0.00
75) Chrysene-d12	14.46	240	301307	20.00	ng	0.00
86) Perylene-d12	16.08	264	259205	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.33	112	883130	110.31	ng	0.02
7) Phenol-d6	5.40	99	1130719	119.80	ng	0.00
23) Nitrobenzene-d5	6.42	82	695910	85.08	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	216111	134.67	ng	0.00
44) 2-Fluorobiphenyl	8.59	172	1202714	86.25	ng	0.00
78) Terphenyl-d14	13.20	244	962441	85.38	ng	0.00
Target Compounds						Qvalue
10) Phenol	5.41	94	27850	2.40	ng	98
49) Dimethylphthalate	9.08	163	351490	22.99	ng	99
70) Phenanthrene	11.24	178	211927	9.46	ng	97
71) Anthracene	11.31	178	56609	2.44	ng	97
74) Fluoranthene	12.70	202	236081	10.17	ng	97
77) Pyrene	12.98	202	212851	10.11	ng	98
80) Benzo(a)anthracene	14.45	228	83890	4.79	ng	98
82) Chrysene	14.49	228	84306	5.27	ng	95
85) Indeno(1,2,3-cd)pyrene	17.35	276	43539	2.86	ng	94
87) Benzo(b)fluoranthene	15.68	252	87767m	5.54	ng	
89) Benzo(a)pyrene	16.02	252	65052	4.41	ng	98
91) Benzo(g,h,i)perylene	17.72	276	39480	3.17	ng	98

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

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