

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096499.D
 Acq On : 5 Jul 2017 23:12
 Operator : SJ/MA
 Sample : I3998-01DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Q5-NSWDL

Quant Time: Jul 06 03:02:08 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 05 13:27:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	98600	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	319167	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	104604	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	155665	20.00	ng	0.00
75) Chrysene-d12	13.86	240	177032	20.00	ng	0.00
86) Perylene-d12	15.27	264	165773	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	120240	18.00	ng	-0.01
7) Phenol-d6	6.34	99	148518	18.08	ng	-0.02
23) Nitrobenzene-d5	7.26	82	81609	15.36	ng	-0.02
41) 2,4,6-Tribromophenol	10.53	330	13906	17.02	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	114848	18.77	ng	0.00
78) Terphenyl-d14	12.81	244	71233	8.60	ng	0.00
Target Compounds						
71) Anthracene	11.31	178	20193	2.35	ng	Qvalue # 84
74) Fluoranthene	12.45	202	181823	22.04	ng	96
77) Pyrene	12.67	202	179377	12.62	ng	99
80) Benzo(a)anthracene	13.85	228	69682	6.80	ng	99
82) Chrysene	13.89	228	57630	5.37	ng	95
85) Indeno(1,2,3-cd)pyrene	16.62	276	22601	2.67	ng	91
87) Benzo(b)fluoranthene	14.87	252	67440m	6.49	ng	
88) Benzo(k)fluoranthene	14.89	252	21692m	2.33	ng	
89) Benzo(a)pyrene	15.21	252	45319	4.89	ng	# 85
91) Benzo(g,h,i)perylene	17.03	276	19257	2.55	ng	94

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

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