

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096516.D
 Acq On : 6 Jul 2017 7:01
 Operator : SJ/MA
 Sample : I3975-08 2X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B8-0-2

Quant Time: Jul 06 15:04:59 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	102471	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	367875	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	133293	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	212961	20.00	ng	0.00
75) Chrysene-d12	13.86	240	188637	20.00	ng	0.00
86) Perylene-d12	15.27	264	122752	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	386827	55.72	ng	0.00
7) Phenol-d6	6.35	99	449202	52.63	ng	-0.01
23) Nitrobenzene-d5	7.27	82	233872	38.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	57225	54.96	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	419572	53.82	ng	-0.01
78) Terphenyl-d14	12.82	244	324562	36.77	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	179095	18.00	ng	98
70) Phenanthrene	11.25	178	51660	4.49	ng	99
74) Fluoranthene	12.44	202	63860	5.66	ng	97
77) Pyrene	12.66	202	60881	4.02	ng	98
82) Chrysene	13.89	228	26242m	2.29	ng	
87) Benzo(b)fluoranthene	14.87	252	22378m	2.91	ng	

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

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