

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040517\
 Data File : BF094216.D
 Acq On : 6 Apr 2017 4:36
 Operator : SJ/MA
 Sample : I2536-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-BB8-BASE

Manual Integrations
 APPROVED

mohammad
 4/6/2017 3:38:51 PM

Quant Time: Apr 06 07:31:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 14:38:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.87	152	154899	20.00	ng	-0.04
21) Naphthalene-d8	7.37	136	541848	20.00	ng	-0.04
38) Acenaphthene-d10	9.52	164	206312	20.00	ng	-0.04
63) Phenanthrene-d10	11.34	188	343456	20.00	ng	-0.04
75) Chrysene-d12	14.59	240	315959	20.00	ng	-0.04
86) Perylene-d12	16.21	264	223995	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	4.40	112	851055	92.80	ng	-0.04
7) Phenol-d6	5.47	99	1060868	93.51	ng	-0.04
23) Nitrobenzene-d5	6.52	82	633783	66.75	ng	-0.04
41) 2,4,6-Tribromophenol	10.49	330	156878	96.23	ng	-0.03
44) 2-Fluorobiphenyl	8.70	172	1055095	78.00	ng	-0.04
78) Terphenyl-d14	13.32	244	745604	52.89	ng	-0.04
Target Compounds						
10) Phenol	5.48	94	114353	8.86	ng	Qvalue 97
17) 2-Methylphenol	6.17	107	159910	18.52	ng	99
20) 3+4-Methylphenols	6.36	107	368597	35.25	ng	# 63
27) 2,4-Dimethylphenol	6.99	122	578576	75.19	ng	98
31) Naphthalene	7.40	128	560173	21.50	ng	# 95
37) 2-Methylnaphthalene	8.25	142	1248084	71.86	ng	98
45) 1,1'-Biphenyl	8.82	154	127249	7.65	ng	98
51) Acenaphthene	9.55	154	43442	3.44	ng	91
54) Dibenzofuran	9.77	168	80935	4.71	ng	# 68
57) Fluorene	10.19	166	88813	6.23	ng	# 82
70) Phenanthrene	11.37	178	276171	14.45	ng	98
71) Anthracene	11.43	178	43068m	2.19	ng	

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

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