

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120318\
 Data File : BG038307.D
 Acq On : 3 Dec 2018 19:41
 Operator : JU/SJ
 Sample : J6060-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20180945-COMPOSITE-4

Quant Time: Dec 04 05:50:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	23927	20.00	ng	-0.02
21) Naphthalene-d8	10.90	136	103638	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	76193	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	195351	20.00	ng	0.00
76) Chrysene-d12	21.75	240	209395	20.00	ng	0.00
87) Perylene-d12	25.06	264	244967	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	174123	125.65	ng	-0.02
7) Phenol-d6	7.22	99	252376	127.58	ng	0.00
23) Nitrobenzene-d5	9.25	82	156680	80.93	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	110496	127.16	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	345568	66.07	ng	-0.02
79) Terphenyl-d14	20.06	244	634808	71.33	ng	0.00
Target Compounds						
10) Phenol	7.25	94	13178	6.289	ng	75
50) Dimethylphthalate	14.16	163	51814	8.578	ng	99
71) Phenanthrene	17.50	178	41867	4.186	ng	99
75) Fluoranthene	19.51	202	59671	5.229	ng	98
78) Pyrene	19.87	202	66537	4.860	ng	97
81) Benzo(a)anthracene	21.72	228	27411	2.166	ng	96
83) Chrysene	21.79	228	28241	2.333	ng	95
88) Benzo(b)fluoranthene	24.00	252	32446	2.407	ng	# 94
90) Benzo(a)pyrene	24.89	252	27573	2.140	ng	# 95

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

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