

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
 Data File : BG038770.D
 Acq On : 20 Dec 2018 20:12
 Operator : JU/SJ
 Sample : J6426-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P001-SS013-3036-01

Quant Time: Dec 21 00:55:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	27464	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	103559	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	68406	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	149279	20.00	ng	0.00
76) Chrysene-d12	21.73	240	157788	20.00	ng	0.00
87) Perylene-d12	25.03	264	173159	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	166471	107.51	ng	0.02
7) Phenol-d6	7.21	99	227045	106.81	ng	0.00
23) Nitrobenzene-d5	9.23	82	145621	71.84	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	109962	116.64	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	347778	69.75	ng	0.00
79) Terphenyl-d14	20.05	244	551381	76.05	ng	0.00

Target Compounds				Qvalue		
10) Phenol	7.24	94	24791	10.977	ng	89
13) 1,4-Dichlorobenzene	8.09	146	10550	4.862	ng	92
14) 1,2-Dichlorobenzene	8.41	146	5799	2.835	ng	94
17) 2-Methylphenol	8.50	107	11137	7.361	ng	94
20) 3+4-Methylphenols	8.82	107	14847	7.105	ng	93
30) 1,2,4-Trichlorobenzene	10.73	180	67989	33.636	ng	96
31) Naphthalene	10.93	128	341900	67.007	ng	99
37) 2-Methylnaphthalene	12.53	142	220611	60.605	ng	98
38) 1-Methylnaphthalene	12.75	142	121972	34.530	ng	99
46) 1,1'-Biphenyl	13.53	154	46555	8.118	ng	96
50) Dimethylphthalate	14.14	163	27651	4.950	ng	# 96
58) Fluorene	15.74	166	18736	3.726	ng	# 95
71) Phenanthrene	17.48	178	97783	12.632	ng	97

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

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