

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100319\
 Data File : BG043012.D
 Acq On : 3 Oct 2019 15:15
 Operator : JU
 Sample : K5168-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20190764-SS-COMPOSITE-4

Quant Time: Oct 04 07:25:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 15:35:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	75941	20.00	ng	-0.01
21) Naphthalene-d8	10.87	136	287912	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	209850	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	469622	20.00	ng	0.00
76) Chrysene-d12	21.69	240	508549	20.00	ng	-0.02
87) Perylene-d12	24.92	264	619655	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	545730	128.25	ng	0.00
7) Phenol-d6	7.24	99	797710	130.18	ng	0.00
23) Nitrobenzene-d5	9.23	82	499217	84.28	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	444774	123.26	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	1173499	81.07	ng	0.00
79) Terphenyl-d14	20.04	244	1951663	81.78	ng	0.00
Target Compounds						
10) Phenol	7.27	94	29995	5.335	ng	87
50) Dimethylphthalate	14.13	163	187564	11.929	ng	99
82) 3,3'-Dichlorobenzidine	21.59	252	36909	2.970	ng	98
86) Indeno(1,2,3-cd)pyrene	28.59	276	107885	2.818	ng	# 92
88) Benzo(b)fluoranthene	23.90	252	71910	2.051	ng	95
91) Dibenzo(a,h)anthracene	28.65	278	88201	2.600	ng	99
92) Benzo(g,h,i)perylene	29.71	276	93941	2.858	ng	# 97

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

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