

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112019\
 Data File : BG043739.D
 Acq On : 20 Nov 2019 22:23
 Operator : JU
 Sample : K5904-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 13-(0-2)

Quant Time: Nov 21 01:20:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 19 15:23:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	24821	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	92774	20.00	ng	0.00
39) Acenaphthene-d10	14.62	164	71013	20.00	ng	0.00
64) Phenanthrene-d10	17.36	188	185882	20.00	ng	0.00
76) Chrysene-d12	21.63	240	193837	20.00	ng	0.00
87) Perylene-d12	24.80	264	225137	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	205252	151.53	ng	0.00
7) Phenol-d6	7.18	99	283901	145.68	ng	0.00
23) Nitrobenzene-d5	9.15	82	164824	91.59	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	165156	122.21	ng	0.00
45) 2-Fluorobiphenyl	13.24	172	425147	82.73	ng	0.00
79) Terphenyl-d14	19.97	244	890525	86.19	ng	0.00
Target Compounds						
10) Phenol	7.21	94	10863	6.100	ng	91
50) Dimethylphthalate	14.07	163	46579	8.468	ng	# 98
71) Phenanthrene	17.41	178	58499	6.146	ng	# 96
75) Fluoranthene	19.42	202	87111	7.020	ng	99
78) Pyrene	19.78	202	73480	6.124	ng	99
81) Benzo(a)anthracene	21.61	228	38929	3.206	ng	94
83) Chrysene	21.67	228	34440	2.855	ng	97
88) Benzo(b)fluoranthene	23.80	252	45866	3.597	ng	92
90) Benzo(a)pyrene	24.66	252	33936	2.787	ng	# 96

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

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