

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082719\
 Data File : BG042523.D
 Acq On : 28 Aug 2019 3:13
 Operator : HP/JU
 Sample : K4526-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T9-12-13-F1-F4-(0-1.9)

Quant Time: Aug 28 04:22:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	43252	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	162092	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	114913	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	257310	20.00	ng	0.00
76) Chrysene-d12	21.68	240	313152	20.00	ng	-0.01
87) Perylene-d12	24.93	264	366193	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	223738	104.77	ng	0.00
7) Phenol-d6	7.16	99	300394	104.13	ng	-0.01
23) Nitrobenzene-d5	9.16	82	175552	74.84	ng	-0.01
42) 2,4,6-Tribromophenol	16.15	330	182918	111.99	ng	-0.01
45) 2-Fluorobiphenyl	13.28	172	500086	73.60	ng	0.00
79) Terphenyl-d14	20.02	244	917754	63.36	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.19	94	12829	4.374	ng	85
50) Dimethylphthalate	14.11	163	37611	4.536	ng	100
75) Fluoranthene	19.47	202	77476	4.967	ng	91
78) Pyrene	19.83	202	74832	3.898	ng	95
81) Benzo(a)anthracene	21.67	228	51878	2.723	ng	98
83) Chrysene	21.73	228	48812	2.657	ng	96
88) Benzo(b)fluoranthene	23.90	252	72430	3.598	ng	# 97
90) Benzo(a)pyrene	24.77	252	51747	2.709	ng	# 96

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

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