

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080119\
 Data File : BG041868.D
 Acq On : 1 Aug 2019 18:09
 Operator : HP/JU
 Sample : K4048-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWL-C4-GW

Manual Integrations
 APPROVED

mohammad
 8/5/2019 1:47:25 PM

Quant Time: Aug 02 03:21:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 08:57:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	39036	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	141868	20.00	ng	-0.01
39) Acenaphthene-d10	14.70	164	93340	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	207272	20.00	ng	0.00
76) Chrysene-d12	21.73	240	245871	20.00	ng	0.00
87) Perylene-d12	25.00	264	278738	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	125784	62.69	ng	0.00
7) Phenol-d6	7.19	99	105793	39.11	ng	0.00
23) Nitrobenzene-d5	9.20	82	209047	101.40	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	159756	150.68	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	573279	108.32	ng	-0.01
79) Terphenyl-d14	20.06	244	1073447	99.90	ng	0.00
Target Compounds						
10) Phenol	7.21	94	32824	11.664	ng	99
32) Benzoic acid	10.10	122	6780m	13.093	ng	
50) Dimethylphthalate	14.15	163	50722	7.957	ng	97
52) Acenaphthene	14.76	154	25358	5.102	ng	96
58) Fluorene	15.75	166	44707	7.328	ng	96
71) Phenanthrene	17.49	178	47558	4.787	ng	94

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

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