

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060319\
 Data File : BG041042.D
 Acq On : 3 Jun 2019 20:03
 Operator : HP/JU
 Sample : K3097-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 0433-HR-13(HACKENSACK)

Manual Integrations
 APPROVED

mohammad
 6/5/2019 8:35:27 AM

Quant Time: Jun 04 09:45:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	67661	20.00	ng	-0.01
21) Naphthalene-d8	10.94	136	277938	20.00	ng	-0.01
39) Acenaphthene-d10	14.75	164	193377	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	453240	20.00	ng	0.00
76) Chrysene-d12	21.77	240	446217	20.00	ng	-0.01
87) Perylene-d12	25.10	264	486461	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	491693	131.04	ng	0.00
7) Phenol-d6	7.27	99	727059	125.46	ng	0.00
23) Nitrobenzene-d5	9.30	82	496300	79.27	ng	0.00
42) 2,4,6-Tribromophenol	16.24	330	356252	124.09	ng	0.00
45) 2-Fluorobiphenyl	13.37	172	880577	65.58	ng	-0.01
79) Terphenyl-d14	20.09	244	1266372	60.23	ng	-0.01
Target Compounds						
10) Phenol	7.29	94	24894	4.007	ng	75
50) Dimethylphthalate	14.19	163	141093	8.619	ng	99
75) Fluoranthene	19.53	202	165739	5.684	ng	99
78) Pyrene	19.90	202	152536	5.259	ng	96
81) Benzo(a)anthracene	21.75	228	143226	4.822	ng	94
83) Chrysene	21.82	228	119615	4.208	ng	99
88) Benzo(b)fluoranthene	24.04	252	154342	5.231	ng	97
90) Benzo(a)pyrene	24.94	252	107811	3.830	ng	97

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

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