

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061421\
 Data File : BG048972.D
 Acq On : 15 Jun 2021 9:13
 Operator : CG/JU
 Sample : M2674-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TMR-DS-01-060921

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:38:49 PM

Quant Time: Jun 15 10:32:37 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 16:03:10 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.120	152	42289	20.00	ng	0.00
21) Naphthalene-d8	10.946	136	162259	20.00	ng	# 0.00
39) Acenaphthene-d10	14.765	164	121010	20.00	ng	0.00
64) Phenanthrene-d10	17.514	188	249511	20.00	ng	# 0.00
76) Chrysene-d12	21.804	240	252727	20.00	ng	#-0.01
86) Perylene-d12	25.147	264	260953	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	340517	125.35	ng	0.00
7) Phenol-d6	7.285	99	418081	102.68	ng	0.02
23) Nitrobenzene-d5	9.295	82	360450	95.79	ng	0.00
42) 2,4,6-Tribromophenol	16.257	330	238545	133.05	ng	0.00
45) 2-Fluorobiphenyl	13.390	172	722294	84.32	ng	0.00
79) Terphenyl-d14	20.123	244	1170817	84.84	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.309	94	178752	42.50	ng	92
17) 2-Methylphenol	8.572	107	243987	88.45	ng	98
20) 3+4-Methylphenols	8.907	107	558261	148.02	ng	93
22) Acetophenone	8.948	105	34105	6.91	ng	# 88
27) 2,4-Dimethylphenol	10.117	122	435041m	166.88	ng	
31) Naphthalene	10.993	128	54589	5.65	ng	98
32) Benzoic acid	10.229	122	18553	10.90	ng	# 77
37) 2-Methylnaphthalene	12.597	142	24519	3.36	ng	93
38) 1-Methylnaphthalene	12.820	142	17425	2.54	ng	# 88
50) Dimethylphthalate	14.212	163	57777	6.13	ng	97
84) Bis(2-ethylhexyl)phtha...	21.698	149	39573	4.01	ng	100

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

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