

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061421\
 Data File : BG048954.D
 Acq On : 14 Jun 2021 18:10
 Operator : CG/JU
 Sample : M2644-08RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1824-25RE

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:38:21 PM

Quant Time: Jun 15 01:13:16 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	8.123	152	44685	20.00	ng	0.00
21) Naphthalene-d8	10.944	136	185225	20.00	ng	# 0.00
39) Acenaphthene-d10	14.763	164	126037	20.00	ng	0.00
64) Phenanthrene-d10	17.512	188	262582	20.00	ng	# 0.00
76) Chrysene-d12	21.801	240	260620	20.00	ng	-0.02
86) Perylene-d12	25.145	264	267238	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.667	112	66178	23.05	ng	0.00
7) Phenol-d6	7.283	99	86835	20.18	ng	0.02
23) Nitrobenzene-d5	9.287	82	167640	39.03	ng	0.00
42) 2,4,6-Tribromophenol	16.249	330	42943	23.00	ng	0.00
45) 2-Fluorobiphenyl	13.388	172	352277	39.48	ng	0.00
79) Terphenyl-d14	20.121	244	530125	37.25	ng	0.00
Target Compounds						
						Qvalue
3) Pyridine	3.958	79	47656m	12.54	ng	
10) Phenol	7.313	94	17866	4.02	ng	81
20) 3+4-Methylphenols	8.881	107	152502	38.27	ng	92
27) 2,4-Dimethylphenol	10.103	122	53797	18.08	ng	92
32) Benzoic acid	10.303	122	72851	37.48	ng	# 73
50) Dimethylphthalate	14.210	163	56750	5.78	ng	# 95

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

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