

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090320\
 Data File : BG046525.D
 Acq On : 3 Sep 2020 20:51
 Operator : CG/JU
 Sample : L3877-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-1

Manual Integrations
 APPROVED

mohammad
 9/4/2020 1:55:03 PM

Quant Time: Sep 04 09:07:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 15:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	83522	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	324942	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	208127	20.00	ng	0.00
64) Phenanthrene-d10	17.30	188	479395	20.00	ng	0.00
76) Chrysene-d12	21.55	240	507803	20.00	ng	0.00
86) Perylene-d12	24.65	264	496713	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	518247	109.90	ng	0.00
7) Phenol-d6	7.11	99	605787	90.62	ng	0.00
23) Nitrobenzene-d5	9.09	82	506010	89.00	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	325474	115.42	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	1197754	87.85	ng	0.00
79) Terphenyl-d14	19.92	244	1859103	77.91	ng	0.00
Target Compounds						
10) Phenol	7.13	94	52575	8.539	ng	94
22) Acetophenone	8.75	105	54935m	6.945	ng	
27) 2,4-Dimethylphenol	9.95	122	14984m	3.676	ng	
31) Naphthalene	10.82	128	7479190	471.406	ng	# 54
37) 2-Methylnaphthalene	12.39	142	1317134	105.262	ng	97
38) 1-Methylnaphthalene	12.61	142	851705m	71.858	ng	
46) 1,1'-Biphenyl	13.39	154	78769	5.440	ng	98

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

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