

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062821\
 Data File : BG049126.D
 Acq On : 28 Jun 2021 19:45
 Operator : CG/JU
 Sample : M2828-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PW-2R

Manual Integrations
 APPROVED

mohammad
 6/29/2021 2:29:24 PM

Quant Time: Jun 29 04:01:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 24 11:02:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.100	152	28018	20.000	ng	# 0.00
21) Naphthalene-d8	10.920	136	113440	20.000	ng	# 0.00
39) Acenaphthene-d10	14.745	164	86928	20.000	ng	0.00
64) Phenanthrene-d10	17.495	188	203190	20.000	ng	# 0.00
76) Chrysene-d12	21.778	240	216181	20.000	ng	-0.01
86) Perylene-d12	25.086	264	232280	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.656	112	101220	56.238	ng	0.00
7) Phenol-d6	7.254	99	84857	31.456	ng	0.00
23) Nitrobenzene-d5	9.269	82	224018	85.152	ng	0.00
42) 2,4,6-Tribromophenol	16.231	330	194773	151.234	ng	0.00
45) 2-Fluorobiphenyl	13.364	172	504327	81.958	ng	0.00
79) Terphenyl-d14	20.097	244	975736	82.661	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	10.180	122	8230m	6.913	ng	
52) Acenaphthene	14.804	154	102208	18.594	ng	93
55) Dibenzofuran	15.144	168	52774	6.176	ng	95
58) Fluorene	15.791	166	138403	19.808	ng	96
71) Phenanthrene	17.536	178	376200	32.905	ng	98
72) Anthracene	17.624	178	42760m	3.722	ng	
75) Fluoranthene	19.539	202	161635	11.724	ng	99
78) Pyrene	19.904	202	109040	6.949	ng	96

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

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