

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062821\
 Data File : BG049129.D
 Acq On : 28 Jun 2021 21:48
 Operator : CG/JU
 Sample : M2828-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ERM-16R

Manual Integrations
 APPROVED

mohammad
 6/29/2021 2:29:30 PM

Quant Time: Jun 29 04:01:52 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	8.103	152	28706	20.000	ng	# 0.00
21) Naphthalene-d8	10.923	136	118764	20.000	ng	# 0.00
39) Acenaphthene-d10	14.743	164	85784	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	196239	20.000	ng	# 0.00
76) Chrysene-d12	21.775	240	217819	20.000	ng	-0.01
86) Perylene-d12	25.089	264	229968	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.653	112	118207	64.102	ng	0.00
7) Phenol-d6	7.251	99	97610	35.316	ng	0.00
23) Nitrobenzene-d5	9.267	82	244505	88.773	ng	0.00
42) 2,4,6-Tribromophenol	16.235	330	212286	167.030	ng	0.00
45) 2-Fluorobiphenyl	13.368	172	538504	88.679	ng	0.00
79) Terphenyl-d14	20.095	244	1041497	87.569	ng	0.00
Target Compounds						
32) Benzoic acid	10.160	122	4698m	3.770	ng	Qvalue
52) Acenaphthene	14.807	154	15684	2.891	ng	# 87
58) Fluorene	15.788	166	23968	3.476	ng	# 84

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

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