

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062821\
 Data File : BG049128.D
 Acq On : 28 Jun 2021 21:07
 Operator : CG/JU
 Sample : M2828-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LT-4

Manual Integrations
 APPROVED

mohammad
 6/29/2021 2:29:27 PM

Quant Time: Jun 29 04:01:42 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	8.104	152	29366	20.000	ng	0.00
21) Naphthalene-d8	10.919	136	126071	20.000	ng	# 0.00
39) Acenaphthene-d10	14.743	164	90166	20.000	ng	# 0.00
64) Phenanthrene-d10	17.493	188	201906	20.000	ng	# 0.00
76) Chrysene-d12	21.776	240	220768	20.000	ng	-0.01
86) Perylene-d12	25.090	264	238618	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.654	112	117262	62.160	ng	0.00
7) Phenol-d6	7.252	99	101163	35.779	ng	0.00
23) Nitrobenzene-d5	9.268	82	244799	83.728	ng	0.00
42) 2,4,6-Tribromophenol	16.230	330	213109	159.529	ng	0.00
45) 2-Fluorobiphenyl	13.369	172	545367	85.445	ng	0.00
79) Terphenyl-d14	20.096	244	1027378	85.228	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	10.172	122	10126m	7.654	ng	
52) Acenaphthene	14.802	154	18821	3.301	ng	91
58) Fluorene	15.789	166	15050	2.077	ng	# 94

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

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