

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010121\
 Data File : BM028324.D
 Acq On : 01 Jan 2021 02:28
 Operator : JU/CG
 Sample : L5139-03DL 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RE115D2-20201215DL

Quant Time: Jan 01 03:00:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-SIM-BM010121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 16:18:03 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.135	152	580	0.40	ng	0.00
7) Naphthalene-d8	10.959	136	2329	0.40	ng	0.00
13) Acenaphthene-d10	14.764	164	1194	0.40	ng	0.00
19) Phenanthrene-d10	17.481	188	2374	0.40	ng	# 0.00
27) Chrysene-d12	21.601	240	2339	0.40	ng	#-0.01
34) Perylene-d12	23.931	264	2240	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.638	112	105	0.05	ng	0.00
5) Phenol-d6	7.273	99	86	0.03	ng	0.00
8) Nitrobenzene-d5	9.299	82	344	0.16	ng	0.00
11) 2-Methylnaphthalene-d10	12.531	152	656	0.16	ng	0.00
14) 2,4,6-Tribromophenol	16.237	330	56	0.09	ng	0.00
15) 2-Fluorobiphenyl	13.390	172	869	0.17	ng	0.00
25) Fluoranthene-d10	19.486	212	1279	0.17	ng	0.00
29) Terphenyl-d14	20.061	244	1178	0.21	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.408	88	1875	2.30	ng	Qvalue 89

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

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