

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061923\
 Data File : BN025904.D
 Acq On : 19 Jun 2023 17:16
 Operator : MA/JU
 Sample : 03317-01DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWOW-BR-05-218-238-061623DL

Quant Time: Jun 19 17:57:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN053123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 14:10:48 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/20/2023
 Supervised By :mohammad ahmed 06/20/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	4734	0.400	ng	0.00
7) Naphthalene-d8	10.825	136	11613	0.400	ng	# 0.01
13) Acenaphthene-d10	14.641	164	6882	0.400	ng	0.01
19) Phenanthrene-d10	17.393	188	14210	0.400	ng	# 0.01
29) Chrysene-d12	21.569	240	9994	0.400	ng	0.00
35) Perylene-d12	24.028	264	11082	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.585	112	112	0.008	ng	0.04
5) Phenol-d6	7.167	99	7	0.000	ng	0.01
8) Nitrobenzene-d5	9.266	82	272m	0.025	ng	0.10
11) 2-Methylnaphthalene-d10	12.460	152	417	0.025	ng	0.06
14) 2,4,6-Tribromophenol	16.152	330	101	0.048	ng	0.02
15) 2-Fluorobiphenyl	13.294	172	778	0.027	ng	0.03
27) Fluoranthene-d10	19.425	212	1449	0.050	ng	0.02
31) Terphenyl-d14	20.006	244	1066	0.044	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.383	88	20280	3.651	ng	# 93

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

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