

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061623\
 Data File : BN025883.D
 Acq On : 16 Jun 2023 14:59
 Operator : MA/JU
 Sample : 03173-13DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 OW-07B-62.5-67.5-061223-FDDL

Quant Time: Jun 16 16:44:31 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 06/19/2023
 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	4046	0.400	ng	0.00
7) Naphthalene-d8	10.825	136	9637	0.400	ng	# 0.01
13) Acenaphthene-d10	14.641	164	5577	0.400	ng	0.01
19) Phenanthrene-d10	17.393	188	10842	0.400	ng	0.01
29) Chrysene-d12	21.569	240	8022	0.400	ng	# 0.00
35) Perylene-d12	24.051	264	8492	0.400	ng	# 0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.571	112	124	0.011	ng	0.03
5) Phenol-d6	7.160	99	3	0.000	ng	0.00
8) Nitrobenzene-d5	9.266	82	174m	0.019	ng	0.10
11) 2-Methylnaphthalene-d10	12.453	152	395	0.028	ng	0.06
14) 2,4,6-Tribromophenol	16.165	330	55	0.032	ng	0.04
15) 2-Fluorobiphenyl	13.304	172	688m	0.029	ng	0.04
27) Fluoranthene-d10	19.426	212	1077	0.048	ng	0.02
31) Terphenyl-d14	20.006	244	871	0.045	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.383	88	14452	3.044	ng	# 93

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

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