

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034539.D
 Acq On : 11 Oct 2024 03:14
 Operator : RC/JU
 Sample : P4241-02
 Misc :
 ALS Vial : 115 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT189D2-GW-20240926

Quant Time: Oct 11 04:05:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 00:21:08 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/11/2024
 Supervised By :mohammad ahmed 10/14/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	3693	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10532	0.400	ng	0.01
13) Acenaphthene-d10	14.030	164	6110	0.400	ng	0.01
19) Phenanthrene-d10	16.783	188	11661	0.400	ng	# 0.00
29) Chrysene-d12	21.008	240	7236	0.400	ng	0.00
35) Perylene-d12	23.113	264	9224	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	939	0.090	ng	0.00
5) Phenol-d6	6.614	99	444m	0.036	ng	0.04
8) Nitrobenzene-d5	8.546	82	1971m	0.245	ng	0.03
11) 2-Methylnaphthalene-d10	11.742	152	4776m	0.307	ng	0.01
14) 2,4,6-Tribromophenol	15.542	330	647	0.234	ng	0.00
15) 2-Fluorobiphenyl	12.651	172	7087	0.285	ng	0.01
27) Fluoranthene-d10	18.829	212	10770	0.366	ng	0.00
31) Terphenyl-d14	19.447	244	8303	0.575	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	25596	5.544	ng	98

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

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