

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101024\
 Data File : BN034547.D
 Acq On : 11 Oct 2024 08:03
 Operator : RC/JU
 Sample : P4231-02DL 2X
 Misc :
 ALS Vial : 123 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D1-20240924DL

Manual Integrations
 APPROVED

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

Quant Time: Oct 11 08:33:54 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.373	152	3856	0.400	ng	0.00
7) Naphthalene-d8	10.137	136	10257	0.400	ng	# 0.01
13) Acenaphthene-d10	14.026	164	6223	0.400	ng	0.01
19) Phenanthrene-d10	16.790	188	11665	0.400	ng	# 0.01
29) Chrysene-d12	21.015	240	8777	0.400	ng	0.00
35) Perylene-d12	23.119	264	9615	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.018	112	642	0.059	ng	0.00
5) Phenol-d6	6.650	99	178m	0.014	ng	0.08
8) Nitrobenzene-d5	8.568	82	1101m	0.140	ng	0.05
11) 2-Methylnaphthalene-d10	11.765	152	2576m	0.170	ng	0.04
14) 2,4,6-Tribromophenol	15.562	330	321	0.114	ng	0.02
15) 2-Fluorobiphenyl	12.657	172	4382m	0.173	ng	0.02
27) Fluoranthene-d10	18.836	212	5955	0.202	ng	0.00
31) Terphenyl-d14	19.449	244	4041	0.231	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.025	88	12697	2.634	ng	99

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

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