

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092324\
 Data File : BN034168.D
 Acq On : 23 Sep 2024 16:07
 Operator : JU/RC
 Sample : P4111-01DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW7-SP100-20240918DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/25/2024
 Supervised By :mohammad ahmed 09/25/2024

Quant Time: Sep 23 16:43:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.423	152	5205	0.400	ng	-0.01
7) Naphthalene-d8	10.180	136	14275	0.400	ng	#-0.02
13) Acenaphthene-d10	14.072	164	8745	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	17906	0.400	ng	0.00
29) Chrysene-d12	21.053	240	11476	0.400	ng	# 0.00
35) Perylene-d12	23.168	264	9028	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.062	112	462	0.031	ng	-0.01
5) Phenol-d6	6.629	99	244	0.014	ng	0.00
8) Nitrobenzene-d5	8.579	82	625	0.057	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	1483	0.070	ng	0.00
14) 2,4,6-Tribromophenol	15.592	330	185	0.047	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	2713m	0.076	ng	0.00
27) Fluoranthene-d10	18.876	212	3976	0.088	ng	-0.01
31) Terphenyl-d14	19.494	244	2887	0.126	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	10098	1.552	ng	# 82

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

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