

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034113.D
 Acq On : 21 Sep 2024 07:30
 Operator : JU/RC
 Sample : P4116-11
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT191D1-20240917

Quant Time: Sep 22 23:56:33 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	7.430	152	5318	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	15902	0.400	ng	#-0.02
13) Acenaphthene-d10	14.072	164	9361	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	21518	0.400	ng	0.00
29) Chrysene-d12	21.053	240	19619	0.400	ng	# 0.00
35) Perylene-d12	23.171	264	19102	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	2172	0.144	ng	0.00
5) Phenol-d6	6.629	99	1696	0.097	ng	0.00
8) Nitrobenzene-d5	8.568	82	3979	0.327	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	7940	0.338	ng	-0.01
14) 2,4,6-Tribromophenol	15.579	330	1271	0.300	ng	0.00
15) 2-Fluorobiphenyl	12.683	172	12864	0.338	ng	-0.02
27) Fluoranthene-d10	18.876	212	22664	0.417	ng	-0.01
31) Terphenyl-d14	19.494	244	15980	0.408	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	11797	1.774	ng	# 81
24) Pentachlorophenol	16.486	266	346	0.077	ng	97

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

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