

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034122.D
 Acq On : 21 Sep 2024 13:38
 Operator : JU/RC
 Sample : P4116-12
 Misc :
 ALS Vial : 86 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT191D2-20240917

Quant Time: Sep 22 23:58:42 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	7.431	152	5888	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	16595	0.400	ng	#-0.02
13) Acenaphthene-d10	14.073	164	9283	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	18544	0.400	ng	0.00
29) Chrysene-d12	21.053	240	12279	0.400	ng	0.00
35) Perylene-d12	23.174	264	10725	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	2106	0.126	ng	0.00
5) Phenol-d6	6.629	99	1424	0.073	ng	0.00
8) Nitrobenzene-d5	8.568	82	3516	0.277	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	7346	0.300	ng	0.00
14) 2,4,6-Tribromophenol	15.580	330	959	0.228	ng	0.00
15) 2-Fluorobiphenyl	12.683	172	11870	0.314	ng	-0.02
27) Fluoranthene-d10	18.876	212	15672	0.335	ng	-0.01
31) Terphenyl-d14	19.494	244	11777	0.480	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	5775	0.785	ng	# 73
24) Pentachlorophenol	16.498	266	131	0.034	ng	88

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

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