

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034124.D
 Acq On : 21 Sep 2024 14:50
 Operator : JU/RC
 Sample : P4116-14
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE125D2-20240917

Quant Time: Sep 22 23:59:07 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	5734	0.400	ng	0.00
7) Naphthalene-d8	10.180	136	16017	0.400	ng	#-0.02
13) Acenaphthene-d10	14.069	164	9577	0.400	ng	-0.01
19) Phenanthrene-d10	16.828	188	20833	0.400	ng	#-0.01
29) Chrysene-d12	21.051	240	14263	0.400	ng	0.00
35) Perylene-d12	23.169	264	11106	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	2092	0.129	ng	0.00
5) Phenol-d6	6.629	99	1186	0.063	ng	0.00
8) Nitrobenzene-d5	8.568	82	3859	0.315	ng	0.00
11) 2-Methylnaphthalene-d10	11.784	152	8495	0.359	ng	0.00
14) 2,4,6-Tribromophenol	15.574	330	1164	0.268	ng	-0.01
15) 2-Fluorobiphenyl	12.690	172	14254	0.366	ng	-0.01
27) Fluoranthene-d10	18.874	212	21079	0.401	ng	-0.01
31) Terphenyl-d14	19.491	244	13907	0.488	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.068	88	63015	8.790	ng	# 88
24) Pentachlorophenol	16.493	266	172	0.040	ng	96

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

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