

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092024\
 Data File : BN034068.D
 Acq On : 20 Sep 2024 03:33
 Operator : JU/RC
 Sample : P3957-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW305I-20240910

Quant Time: Sep 20 04:09:36 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	7.438	152	5432	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	15912	0.400	ng	0.00
13) Acenaphthene-d10	14.086	164	9260	0.400	ng	0.00
19) Phenanthrene-d10	16.840	188	21595	0.400	ng	0.00
29) Chrysene-d12	21.051	240	16421	0.400	ng	0.00
35) Perylene-d12	23.181	264	14571	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	2268	0.147	ng	0.00
5) Phenol-d6	6.636	99	1355	0.076	ng	0.00
8) Nitrobenzene-d5	8.579	82	3673	0.302	ng	0.00
11) 2-Methylnaphthalene-d10	11.791	152	8221	0.350	ng	0.00
14) 2,4,6-Tribromophenol	15.587	330	1330	0.317	ng	0.00
15) 2-Fluorobiphenyl	12.690	172	13373	0.355	ng	-0.01
27) Fluoranthene-d10	18.878	212	24284	0.446	ng	0.00
31) Terphenyl-d14	19.501	244	17359	0.529	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.076	88	6703	0.987	ng	# 81
24) Pentachlorophenol	16.493	266	1037	0.230	ng	99

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

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