

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092024\
 Data File : BN034066.D
 Acq On : 20 Sep 2024 02:21
 Operator : JU/RC
 Sample : P3957-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW203D-20240909

Quant Time: Sep 20 02:45:45 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	6093	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	16816	0.400	ng	0.00
13) Acenaphthene-d10	14.079	164	9167	0.400	ng	0.00
19) Phenanthrene-d10	16.840	188	19291	0.400	ng	0.00
29) Chrysene-d12	21.051	240	12699	0.400	ng	# 0.00
35) Perylene-d12	23.178	264	11247	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	2633	0.152	ng	0.00
5) Phenol-d6	6.629	99	1707	0.085	ng	0.00
8) Nitrobenzene-d5	8.579	82	3999	0.311	ng	0.00
11) 2-Methylnaphthalene-d10	11.791	152	8627	0.347	ng	0.00
14) 2,4,6-Tribromophenol	15.587	330	1355	0.326	ng	0.00
15) 2-Fluorobiphenyl	12.690	172	14205	0.381	ng	-0.01
27) Fluoranthene-d10	18.878	212	20957	0.431	ng	0.00
31) Terphenyl-d14	19.501	244	14060	0.554	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.076	88	1680	0.221	ng	# 36
24) Pentachlorophenol	16.493	266	3092	0.769	ng	99

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

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