

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092024\
 Data File : BN034065.D
 Acq On : 20 Sep 2024 01:44
 Operator : JU/RC
 Sample : P3957-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW202D1-20240909

Quant Time: Sep 20 02:32:23 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	5233	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	15507	0.400	ng	0.00
13) Acenaphthene-d10	14.083	164	8960	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	20931	0.400	ng	# 0.00
29) Chrysene-d12	21.053	240	15933	0.400	ng	0.00
35) Perylene-d12	23.177	264	13544	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	2092	0.141	ng	0.00
5) Phenol-d6	6.636	99	1317	0.076	ng	0.00
8) Nitrobenzene-d5	8.578	82	3368	0.284	ng	0.00
11) 2-Methylnaphthalene-d10	11.791	152	7815	0.341	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	1402	0.345	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	13453	0.369	ng	0.00
27) Fluoranthene-d10	18.880	212	24384	0.462	ng	0.00
31) Terphenyl-d14	19.498	244	16221	0.510	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.075	88	1569	0.240	ng	# 50
24) Pentachlorophenol	16.498	266	495	0.113	ng	97

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

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