

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092124\
 Data File : BN034102.D
 Acq On : 21 Sep 2024 00:51
 Operator : JU/RC
 Sample : P4003-10
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE120D3-20240912

Quant Time: Sep 21 01:21:46 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	7.430	152	4746	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	13169	0.400	ng	0.00
13) Acenaphthene-d10	14.072	164	7922	0.400	ng	0.00
19) Phenanthrene-d10	16.833	188	17737	0.400	ng	0.00
29) Chrysene-d12	21.053	240	13316	0.400	ng	0.00
35) Perylene-d12	23.174	264	11217	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.069	112	1502	0.112	ng	0.00
5) Phenol-d6	6.636	99	1195	0.076	ng	0.00
8) Nitrobenzene-d5	8.578	82	2882	0.286	ng	0.00
11) 2-Methylnaphthalene-d10	11.787	152	6183	0.318	ng	0.00
14) 2,4,6-Tribromophenol	15.579	330	1003	0.279	ng	0.00
15) 2-Fluorobiphenyl	12.693	172	10176	0.316	ng	0.00
27) Fluoranthene-d10	18.876	212	17922	0.400	ng	-0.01
31) Terphenyl-d14	19.498	244	11718	0.441	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.068	88	45067	7.595	ng	Qvalue # 88
24) Pentachlorophenol	16.498	266	270	0.073	ng	97

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

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