

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062223\
 Data File : BN025922.D
 Acq On : 21 Jun 2023 20:58
 Operator : MA/JU
 Sample : SSTD0.409
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4225

Quant Time: Jun 22 01:21:03 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 01:10:16 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.984	152	7888	0.400	ng/u1	0.00
4) Naphthalene-d8	10.798	136	22782	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.621	164	13460	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.364	188	28612	0.400	ng/u1	0.00
17) Chrysene-d12	21.546	240	20683	0.400	ng/u1	0.00
23) Perylene-d12	24.019	264	20858	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.339	96	3609	0.465	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.382	152	14131	0.489	ng/u1	0.00
18) Fluoranthene-d10	19.391	212	30043	0.461	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.373	88	3757	0.453	ng/u1	100
5) Naphthalene	10.847	128	26364	0.455	ng/u1	100
7) 2-Methylnaphthalene	12.453	142	16816	0.482	ng/u1	100
8) 1-Methylnaphthalene	12.673	142	17400	0.454	ng/u1	100
10) Acenaphthylene	14.344	152	25205	0.474	ng/u1	100
11) Acenaphthene	14.681	153	20815	0.469	ng/u1	100
12) Fluorene	15.667	166	23968	0.485	ng/u1	100
14) Pentachlorophenol	17.018	266	4169	1.354	ng/u1	100
15) Phenanthrene	17.406	178	37950	0.455	ng/u1	100
16) Anthracene	17.499	178	33856	0.480	ng/u1	100
19) Fluoranthene	19.419	202	42289	0.447	ng/u1	100
20) Pyrene	19.782	202	44173	0.451	ng/u1	100
21) Benzo(a)anthracene	21.532	228	34184	0.518	ng/u1	100
22) Chrysene	21.584	228	35839	0.475	ng/u1	100
24) Benzo(b)fluoranthene	23.256	252	35305	0.414	ng/u1	100
25) Benzo(k)fluoranthene	23.309	252	35851	0.419	ng/u1	100
26) Benzo(a)pyrene	23.908	252	30654	0.424	ng/u1	100
27) Indeno(1,2,3-cd)pyrene	26.594	276	34355	0.417	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.621	278	25874	0.415	ng/u1	100
29) Benzo(g,h,i)perylene	27.386	276	30546	0.413	ng/u1	100

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

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