

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100225\
 Data File : BN037989.D
 Acq On : 02 Oct 2025 16:23
 Operator : RC/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4272

Quant Time: Oct 02 17:39:20 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 02 09:41:10 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	8750	0.400	ng/u1	0.00
4) Naphthalene-d8	10.612	136	25096	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.465	164	13499	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.214	188	26313	0.400	ng/u1	0.00
17) Chrysene-d12	21.395	240	20590	0.400	ng/u1	0.00
23) Perylene-d12	23.716	264	16984	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	3967	0.427	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.206	152	14092	0.398	ng/u1	0.00
18) Fluoranthene-d10	19.240	212	26000	0.418	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	3956	0.409	ng/u1#	87
5) Naphthalene	10.661	128	27858	0.409	ng/u1	100
7) 2-Methylnaphthalene	12.278	142	18414	0.404	ng/u1	100
8) 1-Methylnaphthalene	12.498	142	19847	0.414	ng/u1	100
10) Acenaphthylene	14.183	152	26093	0.406	ng/u1	99
11) Acenaphthene	14.526	153	19455	0.414	ng/u1	99
12) Fluorene	15.516	166	22096	0.410	ng/u1	100
14) Pentachlorophenol	16.864	266	2163	0.319	ng/u1	99
15) Phenanthrene	17.252	178	35163	0.412	ng/u1	100
16) Anthracene	17.345	178	30321	0.404	ng/u1	100
19) Fluoranthene	19.273	202	36619	0.424	ng/u1	99
20) Pyrene	19.635	202	35622	0.418	ng/u1	98
21) Benzo(a)anthracene	21.380	228	30017	0.402	ng/u1	100
22) Chrysene	21.430	228	32900	0.417	ng/u1	100
24) Benzo(b)fluoranthene	23.014	252	31461	0.425	ng/u1	98
25) Benzo(k)fluoranthene	23.058	252	30688	0.416	ng/u1	99
26) Benzo(a)pyrene	23.613	252	25077	0.410	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	26.082	276	30904	0.420	ng/u1#	92
28) Dibenzo(a,h)anthracene	26.102	278	23823	0.417	ng/u1	99
29) Benzo(g,h,i)perylene	26.806	276	25066	0.412	ng/u1	100

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

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