

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032223\
 Data File : BN024491.D
 Acq On : 22 Mar 2023 22:45
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4251

Quant Time: Mar 23 03:33:26 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 23 03:32:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.029	152	11652	0.400	ng/u1	0.00
4) Naphthalene-d8	10.844	136	34124	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.661	164	18319	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.409	188	37672	0.400	ng/u1	0.00
17) Chrysene-d12	21.594	240	29817	0.400	ng/u1	0.00
23) Perylene-d12	24.099	264	24865	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4379	0.341	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.423	152	16260	0.401	ng/u1	0.00
18) Fluoranthene-d10	19.431	212	32035	0.404	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.430	88	4738	0.340	ng/u1	92
5) Naphthalene	10.894	128	33049	0.388	ng/u1	100
7) 2-Methylnaphthalene	12.500	142	20393	0.395	ng/u1	100
8) 1-Methylnaphthalene	12.714	142	21228	0.391	ng/u1	100
10) Acenaphthylene	14.383	152	29345	0.384	ng/u1	100
11) Acenaphthene	14.726	153	22246	0.378	ng/u1	98
12) Fluorene	15.711	166	24582	0.374	ng/u1	99
14) Pentachlorophenol	17.058	266	3081	0.294	ng/u1#	72
15) Phenanthrene	17.451	178	40065	0.371	ng/u1	99
16) Anthracene	17.540	178	35808	0.378	ng/u1	100
19) Fluoranthene	19.464	202	44955	0.379	ng/u1	98
20) Pyrene	19.822	202	45963	0.382	ng/u1	99
21) Benzo(a)anthracene	21.577	228	37376	0.384	ng/u1	100
22) Chrysene	21.632	228	37908	0.369	ng/u1	100
24) Benzo(b)fluoranthene	23.330	252	36839	0.388	ng/u1	96
25) Benzo(k)fluoranthene	23.380	252	35094	0.356	ng/u1	97
26) Benzo(a)pyrene	23.977	252	30124	0.369	ng/u1	93
27) Indeno(1,2,3-cd)pyrene	26.706	276	34145	0.371	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.730	278	26260	0.375	ng/u1	97
29) Benzo(g,h,i)perylene	27.514	276	29477	0.369	ng/u1	97

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

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