

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103122\
 Data File : BN022442.D
 Acq On : 01 Nov 2022 17:55
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4334

Quant Time: Nov 02 00:55:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 01 01:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.927	152	4304	0.400	ng/u1	0.00
4) Naphthalene-d8	10.750	136	14448	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.600	164	7532	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.357	188	15337	0.400	ng/u1	0.00
17) Chrysene-d12	21.555	240	8846	0.400	ng/u1	0.00
23) Perylene-d12	24.005	264	6411	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	2171	0.388	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.351	152	8524	0.390	ng/u1	0.00
18) Fluoranthene-d10	19.389	212	13638	0.459	ng/u1	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.278	88	2286	0.398	ng/u1	92
5) Naphthalene	10.805	128	16568	0.388	ng/u1	100
7) 2-Methylnaphthalene	12.422	142	10440	0.388	ng/u1	100
8) 1-Methylnaphthalene	12.642	142	9610	0.347	ng/u1	100
10) Acenaphthylene	14.322	152	12358	0.365	ng/u1	100
11) Acenaphthene	14.660	153	10847	0.380	ng/u1	99
12) Fluorene	15.650	166	12719	0.384	ng/u1	98
14) Pentachlorophenol	17.011	266	1539	0.422	ng/u1	100
15) Phenanthrene	17.395	178	19326	0.375	ng/u1	99
16) Anthracene	17.488	178	16359	0.377	ng/u1	99
19) Fluoranthene	19.421	202	18399	0.454	ng/u1	99
20) Pyrene	19.784	202	18158	0.446	ng/u1	98
21) Benzo(a)anthracene	21.541	228	12715	0.384	ng/u1	98
22) Chrysene	21.596	228	13302	0.379	ng/u1	99
24) Benzo(b)fluoranthene	23.254	252	11655	0.362	ng/u1	94
25) Benzo(k)fluoranthene	23.303	252	11537	0.369	ng/u1	95
26) Benzo(a)pyrene	23.894	252	9871	0.385	ng/u1#	92
27) Indeno(1,2,3-cd)pyrene	26.568	276	12594	0.395	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.591	278	9930	0.390	ng/u1	95
29) Benzo(g,h,i)perylene	27.359	276	10533	0.405	ng/u1	98

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

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