

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080822\
 Data File : BN021080.D
 Acq On : 09 Aug 2022 01:54
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4275

Quant Time: Aug 09 02:56:47 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 23:08:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.801	152	2951	0.400	ng/u1	0.01
4) Naphthalene-d8	10.596	136	10024	0.400	ng/u1	0.01
9) Acenaphthene-d10	14.452	164	5581	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.206	188	10661	0.400	ng/u1	0.00
17) Chrysene-d12	21.407	240	8172	0.400	ng/u1	-0.01
23) Perylene-d12	23.748	264	6311	0.400	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1408	0.373	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.202	152	6034	0.411	ng/u1	0.01
18) Fluoranthene-d10	19.245	212	11157	0.442	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.245	88	1482	0.381	ng/u1	96
5) Naphthalene	10.646	128	10725	0.348	ng/u1	99
7) 2-Methylnaphthalene	12.274	142	6640	0.355	ng/u1	99
8) 1-Methylnaphthalene	12.488	142	7257	0.388	ng/u1	99
10) Acenaphthylene	14.170	152	9231	0.354	ng/u1	99
11) Acenaphthene	14.517	153	7434	0.344	ng/u1	98
12) Fluorene	15.507	166	8106	0.344	ng/u1	98
14) Pentachlorophenol	16.876	266	614	0.373	ng/u1	99
15) Phenanthrene	17.248	178	12786	0.335	ng/u1	100
16) Anthracene	17.345	178	10517	0.356	ng/u1	99
19) Fluoranthene	19.273	202	13526	0.374	ng/u1	100
20) Pyrene	19.635	202	13892	0.372	ng/u1	97
21) Benzo(a)anthracene	21.395	228	9498	0.356	ng/u1	99
22) Chrysene	21.445	228	11351	0.323	ng/u1	99
24) Benzo(b)fluoranthene	23.038	252	10012	0.343	ng/u1	98
25) Benzo(k)fluoranthene	23.084	252	10200	0.331	ng/u1	99
26) Benzo(a)pyrene	23.646	252	9099	0.341	ng/u1#	95
27) Indeno(1,2,3-cd)pyrene	26.172	276	9546	0.300	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.186	278	7636	0.304	ng/u1	97
29) Benzo(g,h,i)perylene	26.907	276	8344	0.294	ng/u1	98

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

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