

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082222\
 Data File : BN021321.D
 Acq On : 23 Aug 2022 05:21
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4295

Quant Time: Aug 23 06:04:38 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 18:06:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	3062	0.400	ng/ul	0.00
4) Naphthalene-d8	10.927	136	9829	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.749	164	5353	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.498	188	11294	0.400	ng/ul	0.00
17) Chrysene-d12	21.685	240	9616	0.400	ng/ul	# 0.00
23) Perylene-d12	24.217	264	8597	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.385	96	1486	0.376	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.511	152	5886	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.525	212	11680	0.394	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.422	88	1447	0.350	ng/ul#	90
5) Naphthalene	10.977	128	10723	0.345	ng/ul	98
7) 2-Methylnaphthalene	12.588	142	6788	0.349	ng/ul	98
8) 1-Methylnaphthalene	12.803	142	7254	0.397	ng/ul	100
10) Acenaphthylene	14.472	152	9206	0.333	ng/ul	98
11) Acenaphthene	14.809	153	7180	0.339	ng/ul	98
12) Fluorene	15.795	166	8020	0.339	ng/ul	99
14) Pentachlorophenol	17.143	266	639	0.512	ng/ul	99
15) Phenanthrene	17.540	178	13900	0.332	ng/ul	99
16) Anthracene	17.633	178	11058	0.330	ng/ul	99
19) Fluoranthene	19.552	202	15376	0.337	ng/ul	98
20) Pyrene	19.915	202	14935	0.327	ng/ul	95
21) Benzo(a)anthracene	21.668	228	11573	0.347	ng/ul	100
22) Chrysene	21.726	228	14470	0.338	ng/ul	99
24) Benzo(b)fluoranthene	23.439	252	13577	0.368	ng/ul	97
25) Benzo(k)fluoranthene	23.489	252	13960	0.351	ng/ul	97
26) Benzo(a)pyrene	24.103	252	11799	0.339	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.888	276	14755	0.361	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.921	278	12616	0.376	ng/ul	94
29) Benzo(g,h,i)perylene	27.716	276	14005	0.345	ng/ul	98

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

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