

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052322\
 Data File : BN019913.D
 Acq On : 23 May 2022 13:32
 Operator : CG/JU
 Sample : SSTD1.641
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6241

Quant Time: May 23 14:38:11 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN052322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 23 13:50:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	3836	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	13172	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	7819	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	15430	0.400	ng/ul	0.00
17) Chrysene-d12	21.400	240	12658	0.400	ng/ul	0.00
23) Perylene-d12	23.738	264	9969	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	5658	1.246	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	34795	1.746	ng/ul	0.00
18) Fluoranthene-d10	19.244	212	71790	1.792	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.353	88	6040	1.253	ng/ul#	75
5) Naphthalene	10.689	128	57105	1.339	ng/ul	98
7) 2-Methylnaphthalene	12.300	142	38417	1.479	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	40774	1.695	ng/ul#	100
10) Acenaphthylene	14.192	152	62655	1.620	ng/ul	99
11) Acenaphthene	14.539	153	41591	1.338	ng/ul	100
12) Fluorene	15.520	166	47776	1.401	ng/ul	99
14) Pentachlorophenol	16.871	266	4442	1.338	ng/ul	89
15) Phenanthrene	17.255	178	76074	1.340	ng/ul	99
16) Anthracene	17.344	178	73463	1.587	ng/ul	98
19) Fluoranthene	19.272	202	87704	1.514	ng/ul	98
20) Pyrene	19.634	202	87035	1.505	ng/ul	98
21) Benzo(a)anthracene	21.385	228	72265	1.528	ng/ul	99
22) Chrysene	21.435	228	72185	1.295	ng/ul	99
24) Benzo(b)fluoranthene	23.031	252	62955	1.245	ng/ul	89
25) Benzo(k)fluoranthene	23.077	252	60256	1.237	ng/ul	95
26) Benzo(a)pyrene	23.636	252	58288	1.353	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	26.144	276	61025	1.135	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.161	278	50109	1.115	ng/ul#	78
29) Benzo(g,h,i)perylene	26.882	276	51841	1.055	ng/ul	98

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

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