

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061323\
 Data File : BN025829.D
 Acq On : 14 Jun 2023 05:24
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4304

Quant Time: Jun 14 13:06:03 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 10 02:01:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	7312	0.400	ng/u1	0.00
4) Naphthalene-d8	10.819	136	21379	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.639	164	10485	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.385	188	16807	0.400	ng/u1	0.00
17) Chrysene-d12	21.566	240	10273	0.400	ng/u1	0.00
23) Perylene-d12	24.048	264	10989	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.356	96	2851	0.397	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.403	152	10155	0.375	ng/u1	0.00
18) Fluoranthene-d10	19.409	212	12863	0.397	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.390	88	3008	0.391	ng/u1	98
5) Naphthalene	10.869	128	20755	0.382	ng/u1	100
7) 2-Methylnaphthalene	12.475	142	12261	0.375	ng/u1	100
8) 1-Methylnaphthalene	12.689	142	13206	0.367	ng/u1	100
10) Acenaphthylene	14.362	152	16820	0.406	ng/u1	100
11) Acenaphthene	14.700	153	13397	0.387	ng/u1	99
12) Fluorene	15.685	166	13761	0.357	ng/u1	99
14) Pentachlorophenol	17.047	266	316	0.175	ng/u1	90
15) Phenanthrene	17.427	178	18999	0.388	ng/u1	99
16) Anthracene	17.520	178	15856	0.383	ng/u1	99
19) Fluoranthene	19.437	202	18524	0.394	ng/u1	100
20) Pyrene	19.800	202	19153	0.394	ng/u1	100
21) Benzo(a)anthracene	21.549	228	13823	0.422	ng/u1	99
22) Chrysene	21.604	228	15415	0.411	ng/u1	99
24) Benzo(b)fluoranthene	23.285	252	16142	0.359	ng/u1	97
25) Benzo(k)fluoranthene	23.338	252	16283	0.361	ng/u1	98
26) Benzo(a)pyrene	23.937	252	14392	0.378	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.651	276	20449	0.471	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.671	278	15282	0.465	ng/u1	97
29) Benzo(g,h,i)perylene	27.436	276	17404	0.447	ng/u1	99

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

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