

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020421.D
 Acq On : 25 Jun 2022 05:59
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4379

Quant Time: Jun 25 06:34:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	3025	0.400	ng/u1	0.00
4) Naphthalene-d8	10.535	136	10461	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.377	164	6876	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	15600	0.400	ng/u1	0.00
17) Chrysene-d12	21.309	240	13881	0.400	ng/u1	0.00
23) Perylene-d12	23.592	264	11597	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	1505	0.429	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.124	152	7311	0.426	ng/u1	0.00
18) Fluoranthene-d10	19.151	212	18833	0.411	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	1549	0.433	ng/u1	98
5) Naphthalene	10.584	128	12003	0.372	ng/u1	99
7) 2-Methylnaphthalene	12.201	142	8175	0.378	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	8704	0.425	ng/u1	99
10) Acenaphthylene	14.095	152	13009	0.399	ng/u1	100
11) Acenaphthene	14.442	153	9780	0.375	ng/u1	98
12) Fluorene	15.427	166	11670	0.377	ng/u1	100
14) Pentachlorophenol	16.786	266	1073	0.341	ng/u1	99
15) Phenanthrene	17.162	178	19969	0.368	ng/u1	99
16) Anthracene	17.251	178	18066	0.386	ng/u1	100
19) Fluoranthene	19.178	202	22901	0.361	ng/u1	100
20) Pyrene	19.541	202	23827	0.362	ng/u1	99
21) Benzo(a)anthracene	21.294	228	19643	0.394	ng/u1	100
22) Chrysene	21.347	228	20640	0.372	ng/u1	100
24) Benzo(b)fluoranthene	22.902	252	19148	0.357	ng/u1	97
25) Benzo(k)fluoranthene	22.949	252	18934	0.349	ng/u1	98
26) Benzo(a)pyrene	23.492	252	18170	0.373	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.926	276	17857	0.368	ng/u1#	98
28) Dibenzo(a,h)anthracene	25.943	278	14502	0.370	ng/u1	99
29) Benzo(g,h,i)perylene	26.637	276	15177	0.364	ng/u1	99

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

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