

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029693.D
 Acq On : 11 Feb 2024 20:44
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4343

Quant Time: Feb 11 23:28:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Feb 11 23:18:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.880	152	5089	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	13904	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	6059	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	11053	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	8088	0.400	ng/ul	0.00
23) Perylene-d12	23.840	264	8076	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	2952	0.465	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	7809	0.441	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	11147	0.419	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.370	88	3207	0.464	ng/ul	99
5) Naphthalene	10.727	128	17302	0.442	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	9837	0.438	ng/ul	100
8) 1-Methylnaphthalene	12.564	142	9910	0.434	ng/ul	99
10) Acenaphthylene	14.243	152	12819	0.434	ng/ul	99
11) Acenaphthene	14.581	153	9879	0.440	ng/ul	98
12) Fluorene	15.571	166	10118	0.431	ng/ul	99
14) Pentachlorophenol	16.917	266	919	0.373	ng/ul#	53
15) Phenanthrene	17.314	178	14758	0.427	ng/ul	100
16) Anthracene	17.407	178	12323	0.424	ng/ul	99
19) Fluoranthene	19.337	202	15100	0.406	ng/ul	98
20) Pyrene	19.699	202	15486	0.407	ng/ul	98
21) Benzo(a)anthracene	21.455	228	12283	0.421	ng/ul	100
22) Chrysene	21.508	228	13517	0.424	ng/ul	99
24) Benzo(b)fluoranthene	23.121	252	13571	0.447	ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	14661	0.435	ng/ul	98
26) Benzo(a)pyrene	23.735	252	12322	0.431	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.289	276	15579	0.449	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.326	278	12105	0.445	ng/ul	97
29) Benzo(g,h,i)perylene	27.043	276	14300	0.454	ng/ul	97

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

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