

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021124\
 Data File : BN029676.D
 Acq On : 11 Feb 2024 09:53
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4342

Quant Time: Feb 11 23:25:53 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	5633	0.400	ng/ul	0.00
4) Naphthalene-d8	10.678	136	15643	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.520	164	6784	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	12261	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	8918	0.400	ng/ul	0.00
23) Perylene-d12	23.843	264	8717	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.336	96	3297	0.469	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	8734	0.439	ng/ul	0.00
18) Fluoranthene-d10	19.304	212	11207	0.382	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.365	88	3468	0.454	ng/ul	98
5) Naphthalene	10.727	128	19343	0.439	ng/ul	99
7) 2-Methylnaphthalene	12.344	142	10932	0.433	ng/ul	99
8) 1-Methylnaphthalene	12.564	142	11180	0.435	ng/ul	99
10) Acenaphthylene	14.243	152	14196	0.429	ng/ul	100
11) Acenaphthene	14.580	153	10914	0.434	ng/ul	99
12) Fluorene	15.575	166	11193	0.425	ng/ul	99
14) Pentachlorophenol	16.917	266	1279	0.468	ng/ul#	48
15) Phenanthrene	17.314	178	16204	0.423	ng/ul	100
16) Anthracene	17.407	178	13891	0.431	ng/ul	99
19) Fluoranthene	19.336	202	16592	0.405	ng/ul	98
20) Pyrene	19.699	202	16943	0.404	ng/ul	99
21) Benzo(a)anthracene	21.455	228	13139	0.409	ng/ul	100
22) Chrysene	21.508	228	14928	0.424	ng/ul	99
24) Benzo(b)fluoranthene	23.118	252	14693	0.448	ng/ul	97
25) Benzo(k)fluoranthene	23.168	252	16065	0.442	ng/ul	97
26) Benzo(a)pyrene	23.738	252	13356	0.433	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.295	276	16954	0.453	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.322	278	13134	0.447	ng/ul	98
29) Benzo(g,h,i)perylene	27.036	276	15706	0.462	ng/ul	99

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

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