

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035302.D
 Acq On : 26 Nov 2024 05:08
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4351

Quant Time: Nov 26 06:09:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.314	152	2859	0.400	ng/ul	0.00
4) Naphthalene-d8	10.064	136	6576	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.972	164	3523	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.738	188	7252	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.982	240	6091	0.400	ng/ul	# 0.00
23) Perylene-d12	23.077	264	6663	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1135	0.434	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.664	152	3326	0.345	ng/ul	0.00
18) Fluoranthene-d10	18.791	212	6677	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.006	88	1229	0.433	ng/ul	93
5) Naphthalene	10.108	128	6710	0.405	ng/ul	100
7) 2-Methylnaphthalene	11.741	142	4336	0.377	ng/ul	99
8) 1-Methylnaphthalene	11.967	142	4381	0.378	ng/ul	100
10) Acenaphthylene	13.685	152	6174	0.442	ng/ul	99
11) Acenaphthene	14.037	153	4679	0.438	ng/ul	98
12) Fluorene	15.036	166	5149	0.410	ng/ul	98
14) Pentachlorophenol	16.401	266	568	0.336	ng/ul	96
15) Phenanthrene	16.781	178	8230	0.429	ng/ul	99
16) Anthracene	16.874	178	7112	0.400	ng/ul	99
19) Fluoranthene	18.819	202	9205	0.434	ng/ul#	93
20) Pyrene	19.186	202	9420	0.427	ng/ul#	91
21) Benzo(a)anthracene	20.964	228	8388	0.397	ng/ul	99
22) Chrysene	21.017	228	9405	0.442	ng/ul	100
24) Benzo(b)fluoranthene	22.463	252	8979	0.408	ng/ul	95
25) Benzo(k)fluoranthene	22.504	252	10005	0.441	ng/ul#	95
26) Benzo(a)pyrene	22.986	252	8259	0.410	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.121	276	11090	0.408	ng/ul#	86
28) Dibenzo(a,h)anthracene	25.138	278	8401	0.394	ng/ul#	94
29) Benzo(g,h,i)perylene	25.735	276	8824	0.412	ng/ul#	93

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

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