

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026426.D
 Acq On : 09 Jul 2023 16:41
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4346

Quant Time: Jul 09 17:26:52 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	4923	0.400	ng/ul	0.00
4) Naphthalene-d8	10.731	136	17601	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.565	164	9505	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.325	188	14764	0.400	ng/ul	0.00
17) Chrysene-d12	21.513	240	9343	0.400	ng/ul	0.00
23) Perylene-d12	23.966	264	8834	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	2181	0.393	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.326	152	7920	0.293	ng/ul	0.00
18) Fluoranthene-d10	19.353	212	13285	0.374	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	2085	0.354	ng/ul	96
5) Naphthalene	10.781	128	18651	0.370	ng/ul	99
7) 2-Methylnaphthalene	12.398	142	9177	0.286	ng/ul	99
8) 1-Methylnaphthalene	12.612	142	9946	0.299	ng/ul	100
10) Acenaphthylene	14.292	152	16882	0.379	ng/ul	100
11) Acenaphthene	14.630	153	13521	0.371	ng/ul	98
12) Fluorene	15.625	166	13451	0.322	ng/ul	99
14) Pentachlorophenol	16.992	266	1092	0.206	ng/ul	99
15) Phenanthrene	17.368	178	17150	0.352	ng/ul	99
16) Anthracene	17.465	178	14748	0.340	ng/ul	97
19) Fluoranthene	19.386	202	18218	0.365	ng/ul	98
20) Pyrene	19.744	202	19487	0.375	ng/ul	98
21) Benzo(a)anthracene	21.499	228	13334	0.346	ng/ul	99
22) Chrysene	21.554	228	14409	0.359	ng/ul	99
24) Benzo(b)fluoranthene	23.218	252	12164	0.323	ng/ul	93
25) Benzo(k)fluoranthene	23.264	252	13105	0.346	ng/ul	94
26) Benzo(a)pyrene	23.867	252	11643	0.356	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.550	276	15636	0.437	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.563	278	11870	0.438	ng/ul	99
29) Benzo(g,h,i)perylene	27.335	276	14152	0.445	ng/ul	97

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

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