

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102422\
 Data File : BN022289.D
 Acq On : 24 Oct 2022 08:37
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4324

Quant Time: Oct 25 04:04:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:40:14 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	4803	0.400	ng/ul	0.00
4) Naphthalene-d8	10.767	136	16472	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.614	164	9022	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.374	188	16663	0.400	ng/ul	0.00
17) Chrysene-d12	21.576	240	9230	0.400	ng/ul	0.00
23) Perylene-d12	24.034	264	7360	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.253	96	2199	0.352	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.367	152	9787	0.393	ng/ul	0.00
18) Fluoranthene-d10	19.407	212	14716	0.475	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.291	88	2259	0.352	ng/ul	94
5) Naphthalene	10.816	128	18705	0.384	ng/ul	100
7) 2-Methylnaphthalene	12.439	142	11950	0.390	ng/ul	99
8) 1-Methylnaphthalene	12.659	142	12228	0.388	ng/ul	100
10) Acenaphthylene	14.336	152	15620	0.386	ng/ul	100
11) Acenaphthene	14.679	153	12985	0.380	ng/ul	99
12) Fluorene	15.669	166	14149	0.357	ng/ul	100
14) Pentachlorophenol	17.024	266	1174	0.296	ng/ul	100
15) Phenanthrene	17.412	178	21391	0.382	ng/ul	99
16) Anthracene	17.505	178	17770	0.377	ng/ul	99
19) Fluoranthene	19.440	202	19797	0.468	ng/ul	98
20) Pyrene	19.802	202	19433	0.458	ng/ul	100
21) Benzo(a)anthracene	21.559	228	13451	0.389	ng/ul	98
22) Chrysene	21.611	228	14182	0.387	ng/ul	99
24) Benzo(b)fluoranthene	23.277	252	13332	0.361	ng/ul	98
25) Benzo(k)fluoranthene	23.327	252	12934	0.360	ng/ul	98
26) Benzo(a)pyrene	23.923	252	11161	0.379	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.615	276	14690	0.401	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.635	278	11604	0.397	ng/ul	99
29) Benzo(g,h,i)perylene	27.413	276	12415	0.416	ng/ul	99

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

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