

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040623\
 Data File : BN024802.D
 Acq On : 07 Apr 2023 05:11
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4261

Quant Time: Apr 07 05:40:27 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 06 22:57:45 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	9068	0.400	ng/ul	0.00
4) Naphthalene-d8	10.765	136	27089	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.593	164	14607	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.338	188	32568	0.400	ng/ul	0.00
17) Chrysene-d12	21.529	240	29889	0.400	ng/ul	0.00
23) Perylene-d12	23.987	264	29242	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	3280	0.352	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.354	152	11776	0.352	ng/ul	0.00
18) Fluoranthene-d10	19.368	212	24808	0.339	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.356	88	3627	0.370	ng/ul	96
5) Naphthalene	10.814	128	23639	0.355	ng/ul	100
7) 2-Methylnaphthalene	12.425	142	14388	0.347	ng/ul	100
8) 1-Methylnaphthalene	12.640	142	15249	0.352	ng/ul	100
10) Acenaphthylene	14.316	152	18074	0.350	ng/ul	100
11) Acenaphthene	14.658	153	16304	0.364	ng/ul	100
12) Fluorene	15.644	166	18462	0.361	ng/ul	98
14) Pentachlorophenol	16.988	266	7308	1.038	ng/ul	98
15) Phenanthrene	17.381	178	30810	0.345	ng/ul	100
16) Anthracene	17.474	178	25427	0.341	ng/ul	100
19) Fluoranthene	19.396	202	34537	0.328	ng/ul	97
20) Pyrene	19.758	202	35893	0.335	ng/ul	97
21) Benzo(a)anthracene	21.511	228	33291	0.371	ng/ul	100
22) Chrysene	21.567	228	35701	0.361	ng/ul	99
24) Benzo(b)fluoranthene	23.233	252	40727	0.335	ng/ul	92
25) Benzo(k)fluoranthene	23.286	252	38195	0.327	ng/ul	99
26) Benzo(a)pyrene	23.873	252	34819	0.364	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.537	276	41766	0.367	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.561	278	32666	0.359	ng/ul	98
29) Benzo(g,h,i)perylene	27.329	276	32789	0.335	ng/ul	99

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

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