

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041323\
 Data File : BN024863.D
 Acq On : 13 Apr 2023 16:33
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4268

Quant Time: Apr 14 02:20:42 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 02:44:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	7884	0.400	ng/u1	0.00
4) Naphthalene-d8	10.759	136	21332	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.593	164	11823	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.338	188	26455	0.400	ng/u1	0.00
17) Chrysene-d12	21.528	240	22738	0.400	ng/u1	0.00
23) Perylene-d12	23.975	264	19964	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	3895	0.452	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.348	152	12072	0.441	ng/u1	0.00
18) Fluoranthene-d10	19.367	212	25888	0.458	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.347	88	4222	0.445	ng/u1	98
5) Naphthalene	10.808	128	23957	0.441	ng/u1	99
7) 2-Methylnaphthalene	12.420	142	14838	0.439	ng/u1	100
8) 1-Methylnaphthalene	12.640	142	15605	0.435	ng/u1	99
10) Acenaphthylene	14.311	152	19352	0.436	ng/u1	99
11) Acenaphthene	14.653	153	16765	0.441	ng/u1	99
12) Fluorene	15.639	166	19065	0.448	ng/u1	99
14) Pentachlorophenol	16.992	266	2366	0.419	ng/u1	99
15) Phenanthrene	17.380	178	31746	0.426	ng/u1	100
16) Anthracene	17.473	178	26418	0.426	ng/u1	99
19) Fluoranthene	19.400	202	35755	0.457	ng/u1	99
20) Pyrene	19.762	202	36512	0.455	ng/u1	99
21) Benzo(a)anthracene	21.514	228	31791	0.443	ng/u1	99
22) Chrysene	21.566	228	34679	0.446	ng/u1	100
24) Benzo(b)fluoranthene	23.224	252	34006	0.425	ng/u1	100
25) Benzo(k)fluoranthene	23.273	252	33848	0.433	ng/u1	99
26) Benzo(a)pyrene	23.864	252	27160	0.409	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	26.513	276	31797	0.406	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.533	278	24831	0.403	ng/u1	100
29) Benzo(g,h,i)perylene	27.295	276	28470	0.430	ng/u1	99

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

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