

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013123\
 Data File : BN023884.D
 Acq On : 01 Feb 2023 06:17
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4375

Quant Time: Feb 01 06:57:05 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 01:00:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.859	152	5399	0.400	ng/ul	0.00
4) Naphthalene-d8	10.667	136	14211	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.507	164	8616	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.256	188	19889	0.400	ng/ul	0.00
17) Chrysene-d12	21.450	240	18174	0.400	ng/ul	0.00
23) Perylene-d12	23.850	264	14850	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.243	96	2011	0.347	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	7365	0.331	ng/ul	0.00
18) Fluoranthene-d10	19.286	212	18408	0.377	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.281	88	2123	0.346	ng/ul#	88
5) Naphthalene	10.716	128	13470	0.347	ng/ul	99
7) 2-Methylnaphthalene	12.327	142	8604	0.340	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	8876	0.343	ng/ul	99
10) Acenaphthylene	14.230	152	11337	0.344	ng/ul	99
11) Acenaphthene	14.572	153	10012	0.345	ng/ul	97
12) Fluorene	15.557	166	11816	0.337	ng/ul	98
14) Pentachlorophenol	16.905	266	2035	0.310	ng/ul	99
15) Phenanthrene	17.298	178	20046	0.338	ng/ul	100
16) Anthracene	17.387	178	16754	0.338	ng/ul	99
19) Fluoranthene	19.314	202	23962	0.397	ng/ul	96
20) Pyrene	19.677	202	24402	0.390	ng/ul#	96
21) Benzo(a)anthracene	21.433	228	22317	0.361	ng/ul	100
22) Chrysene	21.485	228	23526	0.362	ng/ul	100
24) Benzo(b)fluoranthene	23.119	252	23798	0.363	ng/ul	98
25) Benzo(k)fluoranthene	23.163	252	23658	0.369	ng/ul	99
26) Benzo(a)pyrene	23.739	252	19465	0.356	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	26.326	276	24634	0.321	ng/ul	97
28) Dibenzo(a,h)anthracene	26.346	278	19272	0.314	ng/ul	98
29) Benzo(g,h,i)perylene	27.094	276	20444	0.324	ng/ul	99

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

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