

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025029.D
 Acq On : 20 Apr 2023 10:08
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4273

Quant Time: Apr 21 04:13:11 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 21 04:13:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	6209	0.400	ng/u1	0.00
4) Naphthalene-d8	10.743	136	19244	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.575	164	11031	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.322	188	24329	0.400	ng/u1	0.00
17) Chrysene-d12	21.515	240	18880	0.400	ng/u1	0.00
23) Perylene-d12	23.964	264	15576	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	3016	0.469	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.332	152	10929	0.442	ng/u1	0.00
18) Fluoranthene-d10	19.350	212	23597	0.474	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.343	88	3105	0.448	ng/u1	98
5) Naphthalene	10.792	128	21470	0.435	ng/u1	100
7) 2-Methylnaphthalene	12.403	142	13323	0.441	ng/u1	100
8) 1-Methylnaphthalene	12.623	142	14209	0.441	ng/u1	100
10) Acenaphthylene	14.293	152	18029	0.439	ng/u1	99
11) Acenaphthene	14.636	153	15539	0.449	ng/u1	99
12) Fluorene	15.626	166	17440	0.450	ng/u1	99
14) Pentachlorophenol	16.976	266	1827	0.383	ng/u1	99
15) Phenanthrene	17.364	178	29627	0.431	ng/u1	99
16) Anthracene	17.457	178	23830	0.424	ng/u1	99
19) Fluoranthene	19.382	202	32189	0.471	ng/u1	99
20) Pyrene	19.745	202	32722	0.466	ng/u1	100
21) Benzo(a)anthracene	21.500	228	25371	0.446	ng/u1	100
22) Chrysene	21.553	228	28511	0.459	ng/u1	100
24) Benzo(b)fluoranthene	23.213	252	26733	0.423	ng/u1	99
25) Benzo(k)fluoranthene	23.262	252	26043	0.428	ng/u1	99
26) Benzo(a)pyrene	23.856	252	21449	0.410	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.511	276	25412	0.422	ng/u1	100
28) Dibenzo(a,h)anthracene	26.538	278	19686	0.418	ng/u1	99
29) Benzo(g,h,i)perylene	27.292	276	22494	0.439	ng/u1	100

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

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