

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062724\
 Data File : BN032301.D
 Acq On : 28 Jun 2024 04:31
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4382

Quant Time: Jun 28 05:07:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 13:23:38 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.620	152	9729	0.400	ng/u1	0.00
4) Naphthalene-d8	10.393	136	29538	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.262	164	15517	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.020	188	28138	0.400	ng/u1	0.00
17) Chrysene-d12	21.220	240	22111	0.400	ng/u1	0.00
23) Perylene-d12	23.439	264	26606	0.400	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.156	96	3899	0.361	ng/u1	0.00
6) 2-Methylnaphthalene-d10	11.993	152	15180	0.343	ng/u1	0.00
18) Fluoranthene-d10	19.055	212	23737	0.357	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.186	88	4337	0.352	ng/u1	94
5) Naphthalene	10.442	128	28524	0.365	ng/u1	100
7) 2-Methylnaphthalene	12.070	142	17780	0.343	ng/u1	99
8) 1-Methylnaphthalene	12.290	142	18282	0.340	ng/u1	99
10) Acenaphthylene	13.985	152	24515	0.380	ng/u1	99
11) Acenaphthene	14.323	153	18122	0.362	ng/u1	99
12) Fluorene	15.322	166	19264	0.323	ng/u1	100
14) Pentachlorophenol	16.682	266	2390	0.375	ng/u1	98
15) Phenanthrene	17.062	178	27615	0.358	ng/u1	100
16) Anthracene	17.155	178	25010	0.363	ng/u1	100
19) Fluoranthene	19.087	202	29854	0.355	ng/u1	100
20) Pyrene	19.450	202	31358	0.356	ng/u1	99
21) Benzo(a)anthracene	21.206	228	29084	0.364	ng/u1	99
22) Chrysene	21.255	228	31525	0.360	ng/u1	98
24) Benzo(b)fluoranthene	22.772	252	33497	0.290	ng/u1	95
25) Benzo(k)fluoranthene	22.816	252	33470	0.298	ng/u1	96
26) Benzo(a)pyrene	23.342	252	28576	0.355	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.676	276	43723	0.347	ng/u1#	99
28) Dibenzo(a,h)anthracene	25.686	278	33698	0.360	ng/u1	97
29) Benzo(g,h,i)perylene	26.361	276	34661	0.346	ng/u1	98

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

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