

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062824\
 Data File : BN032305.D
 Acq On : 28 Jun 2024 09:48
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4383

Quant Time: Jun 28 11:23:43 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	7.616	152	8752	0.400	ng/ul	0.00
4) Naphthalene-d8	10.393	136	27620	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.262	164	15090	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.016	188	25473	0.400	ng/ul	0.00
17) Chrysene-d12	21.220	240	17436	0.400	ng/ul	0.00
23) Perylene-d12	23.438	264	20838	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	3894	0.400	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.988	152	16132	0.389	ng/ul	-0.01
18) Fluoranthene-d10	19.055	212	22133	0.422	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.185	88	4069	0.367	ng/ul	95
5) Naphthalene	10.442	128	30463	0.417	ng/ul	99
7) 2-Methylnaphthalene	12.065	142	19100	0.394	ng/ul	100
8) 1-Methylnaphthalene	12.285	142	20134	0.401	ng/ul	99
10) Acenaphthylene	13.980	152	27606	0.440	ng/ul	99
11) Acenaphthene	14.322	153	20652	0.424	ng/ul	99
12) Fluorene	15.322	166	21194	0.365	ng/ul	99
14) Pentachlorophenol	16.682	266	1965	0.341	ng/ul	99
15) Phenanthrene	17.058	178	28710	0.411	ng/ul	100
16) Anthracene	17.155	178	25456	0.409	ng/ul	100
19) Fluoranthene	19.082	202	28486	0.429	ng/ul	98
20) Pyrene	19.450	202	28819	0.415	ng/ul	99
21) Benzo(a)anthracene	21.206	228	25368	0.403	ng/ul	99
22) Chrysene	21.258	228	27892	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.769	252	28184	0.311	ng/ul	94
25) Benzo(k)fluoranthene	22.816	252	30376	0.345	ng/ul	95
26) Benzo(a)pyrene	23.342	252	28445	0.452	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.676	276	37606	0.381	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.686	278	29378	0.401	ng/ul	97
29) Benzo(g,h,i)perylene	26.357	276	31484	0.401	ng/ul	98

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

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