

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030097.D
 Acq On : 27 Feb 2024 18:20
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4347

Quant Time: Feb 27 23:49:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	3262	0.400	ng/ul	0.00
4) Naphthalene-d8	10.650	136	9286	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	4125	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.263	188	7446	0.400	ng/ul	0.00
17) Chrysene-d12	21.465	240	5654	0.400	ng/ul	0.00
23) Perylene-d12	23.827	264	4858	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	2043	0.502	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	5132	0.434	ng/ul	0.00
18) Fluoranthene-d10	19.294	212	8479	0.456	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.353	88	2237	0.505	ng/ul#	95
5) Naphthalene	10.699	128	11988	0.459	ng/ul	100
7) 2-Methylnaphthalene	12.316	142	6493	0.433	ng/ul	99
8) 1-Methylnaphthalene	12.536	142	6661	0.437	ng/ul	98
10) Acenaphthylene	14.219	152	9013	0.448	ng/ul	98
11) Acenaphthene	14.562	153	7257	0.474	ng/ul	98
12) Fluorene	15.561	166	7444	0.465	ng/ul	98
14) Pentachlorophenol	16.917	266	684	0.412	ng/ul#	47
15) Phenanthrene	17.301	178	10457	0.450	ng/ul	97
16) Anthracene	17.398	178	8725	0.446	ng/ul	98
19) Fluoranthene	19.322	202	11495	0.443	ng/ul	99
20) Pyrene	19.685	202	12312	0.463	ng/ul	98
21) Benzo(a)anthracene	21.451	228	7664	0.376	ng/ul	98
22) Chrysene	21.500	228	10264	0.460	ng/ul	98
24) Benzo(b)fluoranthene	23.108	252	8653	0.473	ng/ul	95
25) Benzo(k)fluoranthene	23.155	252	10776	0.532	ng/ul	97
26) Benzo(a)pyrene	23.725	252	8339	0.485	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.293	276	9481	0.455	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.330	278	7187	0.439	ng/ul#	88
29) Benzo(g,h,i)perylene	27.018	276	9046	0.478	ng/ul	99

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

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