

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018541.D
 Acq On : 02 Feb 2022 09:54
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4317

Quant Time: Feb 02 13:44:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 13:41:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	446	0.400	ng/ul	0.00
4) Naphthalene-d8	10.728	136	1550	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.553	164	836	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.293	188	1666	0.400	ng/ul #	0.00
17) Chrysene-d12	21.467	240	1590	0.400	ng/ul #	0.00
23) Perylene-d12	23.864	264	1409	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.294	96	203	0.395	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.317	152	839	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.318	212	1706	0.384	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.332	88	227	0.420	ng/ul#	82
5) Naphthalene	10.777	128	1679	0.393	ng/ul#	92
7) 2-Methylnaphthalene	12.388	142	1081	0.399	ng/ul	100
8) 1-Methylnaphthalene	12.608	142	1113	0.402	ng/ul	100
10) Acenaphthylene	14.276	152	1284	0.369	ng/ul#	90
11) Acenaphthene	14.618	153	1141	0.395	ng/ul	98
12) Fluorene	15.603	166	1240	0.401	ng/ul#	98
14) Pentachlorophenol	16.947	266	155	1.047	ng/ul	96
15) Phenanthrene	17.336	178	2095	0.390	ng/ul	94
16) Anthracene	17.424	178	1702	0.384	ng/ul#	93
19) Fluoranthene	19.346	202	2367	0.385	ng/ul#	86
20) Pyrene	19.709	202	2418	0.383	ng/ul#	88
21) Benzo(a)anthracene	21.453	228	2076	0.397	ng/ul	96
22) Chrysene	21.502	228	2380	0.395	ng/ul	97
24) Benzo(b)fluoranthene	23.133	252	2277	0.388	ng/ul	90
25) Benzo(k)fluoranthene	23.183	252	2446	0.397	ng/ul#	90
26) Benzo(a)pyrene	23.759	252	1953	0.400	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.356	276	2515	0.395	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.373	278	1973	0.394	ng/ul#	82
29) Benzo(g,h,i)perylene	27.114	276	2184	0.394	ng/ul#	81

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

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