

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020222\
 Data File : BN018567.D
 Acq On : 03 Feb 2022 02:43
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4319

Quant Time: Feb 03 03:56:19 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 14:08:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	449	0.400	ng/ul	0.00
4) Naphthalene-d8	10.728	136	1537	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.554	164	815	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.294	188	1620	0.400	ng/ul #	0.00
17) Chrysene-d12	21.462	240	1483	0.400	ng/ul #	0.00
23) Perylene-d12	23.859	264	1320	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.295	96	204	0.396	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.317	152	825	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.314	212	1654	0.390	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.333	88	246	0.441	ng/ul#	76
5) Naphthalene	10.778	128	1660	0.388	ng/ul#	92
7) 2-Methylnaphthalene	12.389	142	1075	0.394	ng/ul	100
8) 1-Methylnaphthalene	12.609	142	1097	0.392	ng/ul	100
10) Acenaphthylene	14.271	152	1264	0.390	ng/ul#	90
11) Acenaphthene	14.618	153	1118	0.393	ng/ul	99
12) Fluorene	15.599	166	1213	0.394	ng/ul#	97
14) Pentachlorophenol	16.943	266	111	0.368	ng/ul	97
15) Phenanthrene	17.332	178	2047	0.386	ng/ul	94
16) Anthracene	17.425	178	1646	0.387	ng/ul#	92
19) Fluoranthene	19.342	202	2295	0.390	ng/ul#	85
20) Pyrene	19.705	202	2355	0.389	ng/ul#	86
21) Benzo(a)anthracene	21.447	228	1842	0.390	ng/ul	96
22) Chrysene	21.500	228	2246	0.389	ng/ul	97
24) Benzo(b)fluoranthene	23.128	252	2226	0.384	ng/ul	89
25) Benzo(k)fluoranthene	23.172	252	2296	0.378	ng/ul#	90
26) Benzo(a)pyrene	23.751	252	1828	0.390	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.340	276	2409	0.390	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.360	278	1885	0.390	ng/ul#	82
29) Benzo(g,h,i)perylene	27.098	276	2073	0.377	ng/ul#	82

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

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