

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016133.D
 Acq On : 02 Sep 2021 12:03
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4264

Quant Time: Sep 02 13:32:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	2589	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	11253	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	6850	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.251	188	13965	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	14823	0.400	ng/ul	0.00
23) Perylene-d12	23.836	264	14885	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1145	0.380	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	6856	0.432	ng/ul	0.00
18) Fluoranthene-d10	19.281	212	16516	0.415	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	1261	0.383	ng/ul#	91
5) Naphthalene	10.710	128	11410	0.381	ng/ul	99
7) 2-Methylnaphthalene	12.327	142	7940	0.406	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	7974	0.409	ng/ul	100
10) Acenaphthylene	14.219	152	10798	0.417	ng/ul	99
11) Acenaphthene	14.566	153	8252	0.381	ng/ul	99
12) Fluorene	15.552	166	9494	0.382	ng/ul	99
14) Pentachlorophenol	16.904	266	1225	0.542	ng/ul	99
15) Phenanthrene	17.293	178	15222	0.378	ng/ul	99
16) Anthracene	17.381	178	14736	0.424	ng/ul	99
19) Fluoranthene	19.313	202	18870	0.382	ng/ul	98
20) Pyrene	19.676	202	19652	0.393	ng/ul	95
21) Benzo(a)anthracene	21.425	228	19745	0.436	ng/ul	97
22) Chrysene	21.477	228	19522	0.381	ng/ul	97
24) Benzo(b)fluoranthene	23.105	252	20116	0.359	ng/ul	92
25) Benzo(k)fluoranthene	23.155	252	19858	0.352	ng/ul#	93
26) Benzo(a)pyrene	23.731	252	18167	0.382	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.317	276	22598	0.368	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.330	278	18946	0.375	ng/ul#	92
29) Benzo(g,h,i)perylene	27.075	276	19426	0.365	ng/ul	95

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

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