

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016148.D
 Acq On : 02 Sep 2021 22:22
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4265

Quant Time: Sep 03 02:26:41 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	7.867	152	2644	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	11176	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.506	164	6624	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.255	188	13710	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	14451	0.400	ng/ul	0.00
23) Perylene-d12	23.836	264	14325	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1157	0.376	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.261	152	6614	0.420	ng/ul	0.00
18) Fluoranthene-d10	19.285	212	16000	0.413	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	1293	0.384	ng/ul#	89
5) Naphthalene	10.716	128	11374	0.382	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	7816	0.403	ng/ul	98
8) 1-Methylnaphthalene	12.553	142	7817	0.403	ng/ul	99
10) Acenaphthylene	14.224	152	11408	0.455	ng/ul	100
11) Acenaphthene	14.571	153	8252	0.394	ng/ul	98
12) Fluorene	15.556	166	9523	0.397	ng/ul	100
14) Pentachlorophenol	16.908	266	1223	0.551	ng/ul	99
15) Phenanthrene	17.297	178	15317	0.388	ng/ul	99
16) Anthracene	17.386	178	14899	0.437	ng/ul	99
19) Fluoranthene	19.313	202	18936	0.393	ng/ul	100
20) Pyrene	19.675	202	19628	0.403	ng/ul	96
21) Benzo(a)anthracene	21.427	228	19765	0.447	ng/ul	97
22) Chrysene	21.480	228	19445	0.389	ng/ul	97
24) Benzo(b)fluoranthene	23.108	252	20063	0.372	ng/ul	91
25) Benzo(k)fluoranthene	23.155	252	19606	0.362	ng/ul#	91
26) Benzo(a)pyrene	23.731	252	18270	0.400	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.317	276	22454	0.380	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.337	278	18870	0.388	ng/ul#	93
29) Benzo(g,h,i)perylene	27.085	276	19211	0.375	ng/ul	94

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

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