

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015401.D
 Acq On : 08 Jul 2021 09:38
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4110

Quant Time: Jul 08 11:17:38 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.715	152	3010	0.400	ng/ul	0.00
4) Naphthalene-d8	10.485	136	10139	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.339	164	5275	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	10787	0.400	ng/ul	0.00
17) Chrysene-d12	21.287	240	8796	0.400	ng/ul	0.00
23) Perylene-d12	23.528	264	7888	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1354	0.399	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5789	0.364	ng/ul	0.00
18) Fluoranthene-d10	19.122	212	10902	0.377	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	1378	0.376	ng/ul	98
5) Naphthalene	10.529	128	11943	0.369	ng/ul	100
7) 2-Methylnaphthalene	12.146	142	7790	0.368	ng/ul	100
8) 1-Methylnaphthalene	12.371	142	7580	0.370	ng/ul	99
10) Acenaphthylene	14.057	152	9603	0.352	ng/ul	100
11) Acenaphthene	14.400	153	7701	0.368	ng/ul	98
12) Fluorene	15.390	166	8555	0.370	ng/ul	98
14) Pentachlorophenol	16.740	266	858	0.359	ng/ul	99
15) Phenanthrene	17.128	178	14366	0.367	ng/ul	100
16) Anthracene	17.217	178	12179	0.353	ng/ul	100
19) Fluoranthene	19.150	202	15046	0.370	ng/ul	98
20) Pyrene	19.517	202	15197	0.373	ng/ul	99
21) Benzo(a)anthracene	21.272	228	12603	0.387	ng/ul	100
22) Chrysene	21.322	228	14093	0.361	ng/ul	99
24) Benzo(b)fluoranthene	22.856	252	13892	0.377	ng/ul	95
25) Benzo(k)fluoranthene	22.900	252	14466	0.348	ng/ul	99
26) Benzo(a)pyrene	23.429	252	12169	0.393	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.776	276	15435	0.413	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.790	278	12260	0.418	ng/ul	97
29) Benzo(g,h,i)perylene	26.464	276	13148	0.380	ng/ul	99

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

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