

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072121\
 Data File : BN015633.D
 Acq On : 20 Jul 2021 22:02
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4256

Quant Time: Jul 21 01:06:15 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 : Tue Jul 20 18:52:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.779	152	2473	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.567	136	8176	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.418	164	4237	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.170	188	8773	0.400	ng/ul	0.00
17) Chrysene-d12	21.369	240	7385	0.400	ng/ul	# 0.00
23) Perylene-d12	23.710	264	6799	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.277	96	1151	0.413	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.162	152	4824	0.376	ng/ul	0.00
18) Fluoranthene-d10	19.206	212	9674	0.399	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.311	88	1214	0.403	ng/ul#	83
5) Naphthalene	10.611	128	10028	0.384	ng/ul#	91
7) 2-Methylnaphthalene	12.234	142	6395	0.375	ng/ul	99
8) 1-Methylnaphthalene	12.454	142	6367	0.386	ng/ul	99
10) Acenaphthylene	14.136	152	7992	0.365	ng/ul#	90
11) Acenaphthene	14.478	153	6548	0.390	ng/ul	99
12) Fluorene	15.468	166	7217	0.389	ng/ul	99
14) Pentachlorophenol	16.824	266	907	0.467	ng/ul	91
15) Phenanthrene	17.208	178	12232	0.384	ng/ul	97
16) Anthracene	17.301	178	10277	0.366	ng/ul	96
19) Fluoranthene	19.234	202	13245	0.388	ng/ul#	90
20) Pyrene	19.601	202	13621	0.398	ng/ul#	90
21) Benzo(a)anthracene	21.354	228	10949	0.400	ng/ul	99
22) Chrysene	21.404	228	12313	0.376	ng/ul	98
24) Benzo(b)fluoranthene	22.999	252	11837	0.373	ng/ul	93
25) Benzo(k)fluoranthene	23.046	252	13069	0.365	ng/ul#	93
26) Benzo(a)pyrene	23.607	252	10339	0.388	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.118	276	14265	0.443	ng/ul#	50
28) Dibenzo(a,h)anthracene	26.135	278	11319	0.448	ng/ul#	88
29) Benzo(g,h,i)perylene	26.850	276	11480	0.385	ng/ul#	85

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

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