

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015449.D
 Acq On : 09 Jul 2021 16:59
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4114

Quant Time: Jul 09 17:33:50 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	7.711	152	2992	0.400	ng/ul	0.00
4) Naphthalene-d8	10.479	136	10282	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.340	164	5321	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	10490	0.400	ng/ul	0.00
17) Chrysene-d12	21.287	240	8679	0.400	ng/ul	0.00
23) Perylene-d12	23.529	264	8190	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1262	0.374	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	6263	0.388	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	12395	0.435	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	1296	0.356	ng/ul	93
5) Naphthalene	10.529	128	12916	0.394	ng/ul	100
7) 2-Methylnaphthalene	12.146	142	8453	0.394	ng/ul	100
8) 1-Methylnaphthalene	12.366	142	8239	0.397	ng/ul	99
10) Acenaphthylene	14.057	152	11034	0.401	ng/ul	100
11) Acenaphthene	14.400	153	8303	0.394	ng/ul	99
12) Fluorene	15.390	166	9028	0.388	ng/ul	98
14) Pentachlorophenol	16.740	266	1094	0.471	ng/ul	99
15) Phenanthrene	17.128	178	15127	0.397	ng/ul	99
16) Anthracene	17.217	178	13108	0.390	ng/ul	100
19) Fluoranthene	19.150	202	17201	0.429	ng/ul	99
20) Pyrene	19.513	202	16855	0.419	ng/ul	99
21) Benzo(a)anthracene	21.273	228	14150	0.440	ng/ul	99
22) Chrysene	21.322	228	14849	0.386	ng/ul	100
24) Benzo(b)fluoranthene	22.856	252	15134	0.396	ng/ul	94
25) Benzo(k)fluoranthene	22.900	252	14880	0.345	ng/ul	98
26) Benzo(a)pyrene	23.430	252	13426	0.418	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.780	276	16461	0.424	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.790	278	13083	0.430	ng/ul	99
29) Benzo(g,h,i)perylene	26.464	276	13834	0.385	ng/ul	99

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

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