

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016108.D
 Acq On : 27 Aug 2021 16:32
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4262

Quant Time: Aug 27 17:01:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 27 10:53:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	1699	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	7507	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	4712	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	9952	0.400	ng/ul	0.00
17) Chrysene-d12	21.444	240	9761	0.400	ng/ul	0.01
23) Perylene-d12	23.835	264	9070	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	875	0.380	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	4758	0.421	ng/ul	0.00
18) Fluoranthene-d10	19.280	212	10905	0.415	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	840	0.361	ng/ul	98
5) Naphthalene	10.710	128	8267	0.426	ng/ul	100
7) 2-Methylnaphthalene	12.327	142	5762	0.440	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	5320	0.414	ng/ul	99
10) Acenaphthylene	14.219	152	7557	0.454	ng/ul	98
11) Acenaphthene	14.561	153	6094	0.425	ng/ul	99
12) Fluorene	15.547	166	7004	0.412	ng/ul	96
14) Pentachlorophenol	16.895	266	1083	0.520	ng/ul	99
15) Phenanthrene	17.288	178	11527	0.420	ng/ul	99
16) Anthracene	17.377	178	10526	0.444	ng/ul	98
19) Fluoranthene	19.308	202	13504	0.443	ng/ul	99
20) Pyrene	19.670	202	13592	0.443	ng/ul	99
21) Benzo(a)anthracene	21.427	228	13395	0.441	ng/ul	99
22) Chrysene	21.479	228	13686	0.419	ng/ul	100
24) Benzo(b)fluoranthene	23.107	252	13827	0.409	ng/ul	97
25) Benzo(k)fluoranthene	23.154	252	13742	0.403	ng/ul	97
26) Benzo(a)pyrene	23.730	252	12143	0.427	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.306	276	15323	0.407	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.323	278	12769	0.403	ng/ul	98
29) Benzo(g,h,i)perylene	27.064	276	12632	0.389	ng/ul	98

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

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