

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111222\
 Data File : BN022610.D
 Acq On : 11 Nov 2022 13:48
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4339

Quant Time: Nov 11 22:26:41 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 11 22:23:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.898	152	1873	0.400	ng/ul	0.00
4) Naphthalene-d8	10.723	136	5684	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.577	164	3097	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.336	188	6027	0.400	ng/ul	0.00
17) Chrysene-d12	21.544	240	5062	0.400	ng/ul	0.00
23) Perylene-d12	23.985	264	4328	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1088	0.447	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	3577	0.416	ng/ul	0.00
18) Fluoranthene-d10	19.375	212	6732	0.396	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.249	88	1201	0.481	ng/ul#	75
5) Naphthalene	10.772	128	6868	0.409	ng/ul	98
7) 2-Methylnaphthalene	12.400	142	4232	0.400	ng/ul	99
8) 1-Methylnaphthalene	12.615	142	4369	0.401	ng/ul	99
10) Acenaphthylene	14.299	152	5527	0.398	ng/ul	97
11) Acenaphthene	14.637	153	4784	0.408	ng/ul	97
12) Fluorene	15.632	166	5224	0.384	ng/ul	98
14) Pentachlorophenol	16.999	266	584	0.408	ng/ul	99
15) Phenanthrene	17.379	178	8042	0.397	ng/ul	99
16) Anthracene	17.471	178	6834	0.401	ng/ul	97
19) Fluoranthene	19.407	202	8481	0.366	ng/ul	96
20) Pyrene	19.770	202	8675	0.372	ng/ul	99
21) Benzo(a)anthracene	21.529	228	7514	0.396	ng/ul	98
22) Chrysene	21.582	228	7956	0.396	ng/ul	98
24) Benzo(b)fluoranthene	23.236	252	7824	0.360	ng/ul	97
25) Benzo(k)fluoranthene	23.286	252	7866	0.372	ng/ul	96
26) Benzo(a)pyrene	23.877	252	6513	0.376	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.544	276	8937	0.415	ng/ul	92
28) Dibenzo(a,h)anthracene	26.565	278	7161	0.416	ng/ul	94
29) Benzo(g,h,i)perylene	27.329	276	7727	0.440	ng/ul	96

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

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