

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071023\
 Data File : BN026452.D
 Acq On : 10 Jul 2023 18:24
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4349

Quant Time: Jul 11 04:19:52 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3591	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	12876	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.565	164	6682	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.330	188	9736	0.400	ng/ul	0.00
17) Chrysene-d12	21.523	240	6183	0.400	ng/ul	0.00
23) Perylene-d12	23.964	264	6682	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	1686	0.417	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	6408	0.324	ng/ul	0.00
18) Fluoranthene-d10	19.354	212	9221	0.392	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	1659	0.386	ng/ul	94
5) Naphthalene	10.781	128	13845	0.375	ng/ul	98
7) 2-Methylnaphthalene	12.403	142	7376	0.314	ng/ul	99
8) 1-Methylnaphthalene	12.612	142	8340	0.343	ng/ul	100
10) Acenaphthylene	14.288	152	12121	0.387	ng/ul	99
11) Acenaphthene	14.626	153	9939	0.388	ng/ul	99
12) Fluorene	15.625	166	9502	0.324	ng/ul	99
14) Pentachlorophenol	16.996	266	704	0.202	ng/ul	99
15) Phenanthrene	17.368	178	11840	0.368	ng/ul	98
16) Anthracene	17.465	178	9943	0.348	ng/ul	96
19) Fluoranthene	19.387	202	12900	0.391	ng/ul	98
20) Pyrene	19.749	202	13827	0.403	ng/ul	96
21) Benzo(a)anthracene	21.505	228	8503	0.333	ng/ul	98
22) Chrysene	21.558	228	9697	0.365	ng/ul	98
24) Benzo(b)fluoranthene	23.218	252	10237	0.359	ng/ul	91
25) Benzo(k)fluoranthene	23.265	252	11088	0.387	ng/ul	93
26) Benzo(a)pyrene	23.858	252	9300	0.375	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.524	276	11586	0.428	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.534	278	8613	0.421	ng/ul#	92
29) Benzo(g,h,i)perylene	27.299	276	10609	0.441	ng/ul	98

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

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