

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071423\
 Data File : BN026513.D
 Acq On : 14 Jul 2023 23:14
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4357

Quant Time: Jul 15 01:56:55 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 15 01:53:23 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	2782	0.400	ng/ul	0.00
4) Naphthalene-d8	10.759	136	9842	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.585	164	6164	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.335	188	12234	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	10870	0.400	ng/ul	# 0.00
23) Perylene-d12	23.933	264	10406	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	1599	0.436	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.354	152	5580	0.417	ng/ul	0.00
18) Fluoranthene-d10	19.360	212	14151	0.376	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.339	88	1602	0.462	ng/ul#	92
5) Naphthalene	10.809	128	10529	0.390	ng/ul	98
7) 2-Methylnaphthalene	12.425	142	6116	0.399	ng/ul	99
8) 1-Methylnaphthalene	12.640	142	6855	0.417	ng/ul	100
10) Acenaphthylene	14.312	152	11159	0.370	ng/ul	98
11) Acenaphthene	14.650	153	8534	0.371	ng/ul	98
12) Fluorene	15.640	166	9256	0.384	ng/ul	97
14) Pentachlorophenol	16.997	266	1477	0.479	ng/ul	99
15) Phenanthrene	17.377	178	14204	0.365	ng/ul	98
16) Anthracene	17.470	178	12999	0.384	ng/ul	97
19) Fluoranthene	19.387	202	17742	0.348	ng/ul	98
20) Pyrene	19.750	202	19368	0.350	ng/ul	99
21) Benzo(a)anthracene	21.501	228	15076	0.377	ng/ul	98
22) Chrysene	21.553	228	15807	0.348	ng/ul	99
24) Benzo(b)fluoranthene	23.193	252	14101	0.348	ng/ul	96
25) Benzo(k)fluoranthene	23.243	252	15118	0.356	ng/ul	96
26) Benzo(a)pyrene	23.824	252	12891	0.357	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	26.451	276	16786	0.355	ng/ul#	83
28) Dibenzo(a,h)anthracene	26.461	278	13303	0.366	ng/ul#	90
29) Benzo(g,h,i)perylene	27.226	276	14252	0.341	ng/ul#	91

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

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