

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026408.D
 Acq On : 09 Jul 2023 05:04
 Operator : MA/JU
 Sample : PB153851BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS851

Quant Time: Jul 09 05:45:46 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 : Sat Jul 08 09:20:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	4302	0.400	ng/ul	0.00
4) Naphthalene-d8	10.737	136	13647	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.571	164	8453	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.327	188	12759	0.400	ng/ul	0.00
17) Chrysene-d12	21.518	240	8282	0.400	ng/ul	0.00
23) Perylene-d12	23.971	264	8219	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3589	0.741	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	7436	0.354	ng/ul	0.00
18) Fluoranthene-d10	19.355	212	11564	0.367	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	8651	1.682	ng/ul#	86
5) Naphthalene	10.787	128	12242	0.313	ng/ul	99
7) 2-Methylnaphthalene	12.409	142	8810	0.354	ng/ul	99
8) 1-Methylnaphthalene	12.618	142	9816	0.380	ng/ul	99
10) Acenaphthylene	14.293	152	14332	0.362	ng/ul	99
11) Acenaphthene	14.631	153	11641	0.360	ng/ul	99
12) Fluorene	15.626	166	11485	0.309	ng/ul	99
14) Pentachlorophenol	16.993	266	1845	0.403	ng/ul	99
15) Phenanthrene	17.369	178	14874	0.353	ng/ul	98
16) Anthracene	17.466	178	12451	0.332	ng/ul	96
19) Fluoranthene	19.383	202	16038	0.363	ng/ul	99
20) Pyrene	19.750	202	13793	0.300	ng/ul#	90
21) Benzo(a)anthracene	21.501	228	11765	0.344	ng/ul	99
22) Chrysene	21.553	228	13442	0.378	ng/ul	99
24) Benzo(b)fluoranthene	23.219	252	13219	0.377	ng/ul	92
25) Benzo(k)fluoranthene	23.266	252	14083	0.400	ng/ul	93
26) Benzo(a)pyrene	23.868	252	11851	0.389	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.558	276	14591	0.438	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.572	278	11027	0.438	ng/ul	97
29) Benzo(g,h,i)perylene	27.333	276	13034	0.440	ng/ul	97

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

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