

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090523\
 Data File : BN027400.D
 Acq On : 06 Sep 2023 07:49
 Operator : MA/JU
 Sample : PB155237BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS237

Quant Time: Sep 06 08:18:44 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 02 06:16:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	3390	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	11133	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.689	164	6369	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.439	188	15084	0.400	ng/ul	0.00
17) Chrysene-d12	21.615	240	14804	0.400	ng/ul	0.00
23) Perylene-d12	24.141	264	13731	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.418	96	2769	0.614	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	5111	0.302	ng/ul	0.00
18) Fluoranthene-d10	19.455	212	14360	0.283	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.452	88	6697	1.468	ng/ul	89
5) Naphthalene	10.927	128	10229	0.317	ng/ul	98
7) 2-Methylnaphthalene	12.528	142	6138	0.304	ng/ul	100
8) 1-Methylnaphthalene	12.742	142	6590	0.306	ng/ul	99
10) Acenaphthylene	14.416	152	9388	0.322	ng/ul	98
11) Acenaphthene	14.749	153	7870	0.314	ng/ul	99
12) Fluorene	15.739	166	8806	0.313	ng/ul	99
14) Pentachlorophenol	17.067	266	2217	0.754	ng/ul	100
15) Phenanthrene	17.481	178	15095	0.307	ng/ul	99
16) Anthracene	17.574	178	12412	0.318	ng/ul	99
19) Fluoranthene	19.487	202	18884	0.277	ng/ul	100
20) Pyrene	19.850	202	20022	0.289	ng/ul	99
21) Benzo(a)anthracene	21.600	228	16939	0.318	ng/ul	99
22) Chrysene	21.656	228	19277	0.310	ng/ul	99
24) Benzo(b)fluoranthene	23.351	252	18959	0.292	ng/ul	97
25) Benzo(k)fluoranthene	23.407	252	19759	0.312	ng/ul	97
26) Benzo(a)pyrene	24.027	252	15531	0.307	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.817	276	17647	0.314	ng/ul#	45
28) Dibenzo(a,h)anthracene	26.837	278	13514	0.314	ng/ul	98
29) Benzo(g,h,i)perylene	27.636	276	16344	0.322	ng/ul	99

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

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