

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021123.D
 Acq On : 10 Aug 2022 18:24
 Operator : CG/JU
 Sample : PB146872BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS872

Quant Time: Aug 10 22:08:09 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 09 23:10:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	2210	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	7733	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.438	164	4435	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.189	188	8190	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	5892	0.400	ng/ul	0.00
23) Perylene-d12	23.728	264	4331	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2082	0.736	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.186	152	4397	0.388	ng/ul	0.00
18) Fluoranthene-d10	19.231	212	8361	0.460	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	5813	1.996	ng/ul#	85
5) Naphthalene	10.635	128	8169	0.344	ng/ul	100
7) 2-Methylnaphthalene	12.263	142	4929	0.342	ng/ul	100
8) 1-Methylnaphthalene	12.477	142	5425	0.376	ng/ul	99
10) Acenaphthylene	14.161	152	7142	0.344	ng/ul	100
11) Acenaphthene	14.503	153	5787	0.337	ng/ul	99
12) Fluorene	15.498	166	6254	0.334	ng/ul	98
14) Pentachlorophenol	16.859	266	686	0.543	ng/ul	100
15) Phenanthrene	17.231	178	9716	0.332	ng/ul	100
16) Anthracene	17.328	178	7827	0.345	ng/ul	99
19) Fluoranthene	19.259	202	10358	0.397	ng/ul	99
20) Pyrene	19.621	202	10772	0.400	ng/ul	98
21) Benzo(a)anthracene	21.380	228	6825	0.355	ng/ul	100
22) Chrysene	21.433	228	8566	0.338	ng/ul	99
24) Benzo(b)fluoranthene	23.017	252	7071	0.353	ng/ul	96
25) Benzo(k)fluoranthene	23.064	252	7483	0.354	ng/ul	97
26) Benzo(a)pyrene	23.622	252	6517	0.356	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.139	276	6795	0.311	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.156	278	5353	0.310	ng/ul	97
29) Benzo(g,h,i)perylene	26.870	276	6176	0.317	ng/ul	96

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

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