

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042123\
 Data File : BN025059.D
 Acq On : 21 Apr 2023 17:02
 Operator : CG/JU
 Sample : PB152233BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS233

Quant Time: Apr 22 00:17:06 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 21 04:13:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	8857	0.400	ng/u1	0.00
4) Naphthalene-d8	10.737	136	27755	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.571	164	15792	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.322	188	33297	0.400	ng/u1	0.00
17) Chrysene-d12	21.518	240	21467	0.400	ng/u1	0.00
23) Perylene-d12	23.965	264	16954	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	6705	0.731	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.332	152	12816	0.360	ng/u1	0.00
18) Fluoranthene-d10	19.350	212	25536	0.451	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.344	88	16712	1.692	ng/u1	89
5) Naphthalene	10.787	128	24305	0.342	ng/u1	100
7) 2-Methylnaphthalene	12.404	142	15051	0.346	ng/u1	99
8) 1-Methylnaphthalene	12.618	142	15952	0.344	ng/u1	100
10) Acenaphthylene	14.294	152	19935	0.339	ng/u1	100
11) Acenaphthene	14.636	153	17205	0.347	ng/u1	99
12) Fluorene	15.621	166	19000	0.343	ng/u1	99
14) Pentachlorophenol	16.972	266	4079	0.624	ng/u1	99
15) Phenanthrene	17.365	178	31232	0.332	ng/u1	100
16) Anthracene	17.458	178	25281	0.329	ng/u1	100
19) Fluoranthene	19.383	202	33411	0.430	ng/u1	99
20) Pyrene	19.745	202	33657	0.422	ng/u1	99
21) Benzo(a)anthracene	21.504	228	23358	0.361	ng/u1	99
22) Chrysene	21.553	228	26729	0.379	ng/u1	99
24) Benzo(b)fluoranthene	23.216	252	23732	0.345	ng/u1	98
25) Benzo(k)fluoranthene	23.263	252	23654	0.358	ng/u1	98
26) Benzo(a)pyrene	23.860	252	18769	0.330	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.518	276	22193	0.338	ng/u1#	48
28) Dibenzo(a,h)anthracene	26.542	278	17134	0.335	ng/u1	99
29) Benzo(g,h,i)perylene	27.296	276	20193	0.362	ng/u1	99

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

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