

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022891.D
 Acq On : 26 Nov 2022 21:08
 Operator : CG/JU
 Sample : PB149029BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS029

Quant Time: Nov 28 03:49:36 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN112622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 28 03:16:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	7182	0.400	ng/u1	0.00
4) Naphthalene-d8	10.892	136	20042	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.710	164	11275	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.453	188	24490	0.400	ng/u1	0.00
17) Chrysene-d12	21.639	240	16134	0.400	ng/u1	0.00
23) Perylene-d12	24.127	264	10893	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.412	96	5945	0.690	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.476	152	12159	0.398	ng/u1	0.00
18) Fluoranthene-d10	19.480	212	24536	0.412	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.445	88	15415	1.715	ng/u1	92
5) Naphthalene	10.942	128	22468	0.369	ng/u1	100
7) 2-Methylnaphthalene	12.547	142	14241	0.375	ng/u1	100
8) 1-Methylnaphthalene	12.767	142	15158	0.392	ng/u1	99
10) Acenaphthylene	14.432	152	19767	0.356	ng/u1	99
11) Acenaphthene	14.770	153	15871	0.361	ng/u1	99
12) Fluorene	15.755	166	18025	0.362	ng/u1	100
14) Pentachlorophenol	17.099	266	2987	0.583	ng/u1#	100
15) Phenanthrene	17.496	178	30499	0.363	ng/u1	100
16) Anthracene	17.588	178	26370	0.366	ng/u1	99
19) Fluoranthene	19.508	202	31895	0.383	ng/u1	100
20) Pyrene	19.871	202	31536	0.386	ng/u1	100
21) Benzo(a)anthracene	21.622	228	23666	0.384	ng/u1	99
22) Chrysene	21.674	228	25915	0.383	ng/u1	99
24) Benzo(b)fluoranthene	23.364	252	20971	0.380	ng/u1	99
25) Benzo(k)fluoranthene	23.413	252	22495	0.386	ng/u1	99
26) Benzo(a)pyrene	24.016	252	15879	0.378	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.738	276	18253	0.365	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.768	278	14708	0.371	ng/u1	100
29) Benzo(g,h,i)perylene	27.543	276	15844	0.364	ng/u1	99

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

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