

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022893.D
 Acq On : 26 Nov 2022 22:22
 Operator : CG/JU
 Sample : PB149275BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS275

Quant Time: Nov 28 03:49:56 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	8.070	152	7902	0.400	ng/u1	0.00
4) Naphthalene-d8	10.891	136	21005	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.709	164	10612	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.456	188	21174	0.400	ng/u1	0.00
17) Chrysene-d12	21.637	240	13569	0.400	ng/u1	0.00
23) Perylene-d12	24.127	264	10061	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.412	96	5780	0.610	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.475	152	10455	0.327	ng/u1	0.00
18) Fluoranthene-d10	19.479	212	16964	0.338	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.450	88	15152	1.532	ng/u1	94
5) Naphthalene	10.940	128	20837	0.327	ng/u1	100
7) 2-Methylnaphthalene	12.546	142	12590	0.316	ng/u1	99
8) 1-Methylnaphthalene	12.766	142	13409	0.331	ng/u1	99
10) Acenaphthylene	14.431	152	16473	0.315	ng/u1	99
11) Acenaphthene	14.773	153	13189	0.318	ng/u1	100
12) Fluorene	15.754	166	14449	0.308	ng/u1	99
14) Pentachlorophenol	17.097	266	2339	0.528	ng/u1#	100
15) Phenanthrene	17.494	178	23036	0.317	ng/u1	100
16) Anthracene	17.587	178	19295	0.309	ng/u1	99
19) Fluoranthene	19.507	202	22656	0.323	ng/u1	98
20) Pyrene	19.870	202	22399	0.326	ng/u1	99
21) Benzo(a)anthracene	21.619	228	17024	0.328	ng/u1	100
22) Chrysene	21.675	228	19200	0.338	ng/u1	100
24) Benzo(b)fluoranthene	23.364	252	16372	0.321	ng/u1	99
25) Benzo(k)fluoranthene	23.414	252	17697	0.328	ng/u1	99
26) Benzo(a)pyrene	24.013	252	12683	0.327	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.735	276	16543	0.358	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.765	278	13345	0.364	ng/u1	99
29) Benzo(g,h,i)perylene	27.537	276	14740	0.366	ng/u1	100

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

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