

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023482.D
 Acq On : 26 Dec 2022 12:28
 Operator : CG/JU
 Sample : PB149683BS
 Misc :
 ALS Vial : 119 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS683

Quant Time: Dec 27 03:49:23 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 25 05:17:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.985	152	5435	0.400	ng/u1	0.00
4) Naphthalene-d8	10.797	136	17820	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.635	164	9898	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.380	188	20906	0.400	ng/u1	0.00
17) Chrysene-d12	21.569	240	11725	0.400	ng/u1	0.00
23) Perylene-d12	24.013	264	8946	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	3494	0.592	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.387	152	10272	0.373	ng/u1	0.00
18) Fluoranthene-d10	19.410	212	19580	0.462	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.374	88	9485	1.570	ng/u1	91
5) Naphthalene	10.847	128	19170	0.366	ng/u1	100
7) 2-Methylnaphthalene	12.458	142	12176	0.364	ng/u1	100
8) 1-Methylnaphthalene	12.678	142	13039	0.378	ng/u1	100
10) Acenaphthylene	14.352	152	16167	0.369	ng/u1	100
11) Acenaphthene	14.695	153	13475	0.365	ng/u1	99
12) Fluorene	15.680	166	14902	0.359	ng/u1	100
14) Pentachlorophenol	17.030	266	3212	0.554	ng/u1	100
15) Phenanthrene	17.423	178	24329	0.357	ng/u1	100
16) Anthracene	17.511	178	22181	0.393	ng/u1	99
19) Fluoranthene	19.437	202	25601	0.450	ng/u1	100
20) Pyrene	19.800	202	26329	0.459	ng/u1	100
21) Benzo(a)anthracene	21.552	228	18457	0.404	ng/u1	100
22) Chrysene	21.607	228	18830	0.406	ng/u1	100
24) Benzo(b)fluoranthene	23.265	252	16918	0.366	ng/u1	98
25) Benzo(k)fluoranthene	23.312	252	16237	0.374	ng/u1	98
26) Benzo(a)pyrene	23.902	252	13539	0.370	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	26.561	276	17208	0.402	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.587	278	13598	0.410	ng/u1	99
29) Benzo(g,h,i)perylene	27.349	276	14592	0.419	ng/u1	99

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

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