

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022906.D
 Acq On : 27 Nov 2022 06:57
 Operator : CG/JU
 Sample : PB149075BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS075

Quant Time: Nov 28 03:51:29 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	8386	0.400	ng/u1	0.00
4) Naphthalene-d8	10.887	136	23269	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.710	164	12874	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.453	188	27827	0.400	ng/u1	0.00
17) Chrysene-d12	21.633	240	18019	0.400	ng/u1	0.00
23) Perylene-d12	24.124	264	12350	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.412	96	5336	0.530	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.470	152	10246	0.289	ng/u1	0.00
18) Fluoranthene-d10	19.476	212	19712	0.296	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.446	88	13259	1.263	ng/u1	92
5) Naphthalene	10.936	128	19173	0.271	ng/u1	100
7) 2-Methylnaphthalene	12.542	142	12001	0.272	ng/u1	100
8) 1-Methylnaphthalene	12.762	142	12781	0.285	ng/u1	99
10) Acenaphthylene	14.428	152	16044	0.253	ng/u1	99
11) Acenaphthene	14.770	153	13326	0.265	ng/u1	98
12) Fluorene	15.756	166	15026	0.264	ng/u1	99
14) Pentachlorophenol	17.099	266	2062	0.354	ng/u1#	100
15) Phenanthrene	17.496	178	25251	0.264	ng/u1	100
16) Anthracene	17.584	178	21386	0.261	ng/u1	100
19) Fluoranthene	19.504	202	25555	0.274	ng/u1	98
20) Pyrene	19.866	202	25109	0.275	ng/u1	98
21) Benzo(a)anthracene	21.619	228	18191	0.264	ng/u1	99
22) Chrysene	21.671	228	20865	0.276	ng/u1	99
24) Benzo(b)fluoranthene	23.358	252	16601	0.265	ng/u1	97
25) Benzo(k)fluoranthene	23.408	252	18322	0.277	ng/u1	98
26) Benzo(a)pyrene	24.010	252	12231	0.257	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.731	276	14524	0.256	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.755	278	11870	0.264	ng/u1	99
29) Benzo(g,h,i)perylene	27.536	276	12595	0.255	ng/u1	99

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

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