

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022904.D
 Acq On : 27 Nov 2022 05:44
 Operator : CG/JU
 Sample : PB149073BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS073

Quant Time: Nov 28 03:51:12 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	7222	0.400	ng/u1	0.00
4) Naphthalene-d8	10.887	136	19972	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.710	164	10952	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.453	188	23875	0.400	ng/u1	0.00
17) Chrysene-d12	21.633	240	15865	0.400	ng/u1	0.00
23) Perylene-d12	24.121	264	10298	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.408	96	5328	0.615	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.470	152	10515	0.345	ng/u1	0.00
18) Fluoranthene-d10	19.476	212	20526	0.350	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.446	88	13688	1.514	ng/u1	92
5) Naphthalene	10.936	128	20186	0.333	ng/u1	100
7) 2-Methylnaphthalene	12.542	142	12590	0.332	ng/u1	100
8) 1-Methylnaphthalene	12.762	142	13483	0.350	ng/u1	100
10) Acenaphthylene	14.428	152	17155	0.318	ng/u1	100
11) Acenaphthene	14.770	153	14095	0.330	ng/u1	99
12) Fluorene	15.756	166	16055	0.332	ng/u1	99
14) Pentachlorophenol	17.095	266	2322	0.465	ng/u1#	100
15) Phenanthrene	17.496	178	26938	0.329	ng/u1	99
16) Anthracene	17.584	178	22538	0.320	ng/u1	99
19) Fluoranthene	19.504	202	27843	0.340	ng/u1	99
20) Pyrene	19.866	202	27216	0.339	ng/u1	98
21) Benzo(a)anthracene	21.616	228	20276	0.334	ng/u1	100
22) Chrysene	21.671	228	22934	0.345	ng/u1	100
24) Benzo(b)fluoranthene	23.361	252	18348	0.352	ng/u1	98
25) Benzo(k)fluoranthene	23.411	252	19334	0.351	ng/u1	99
26) Benzo(a)pyrene	24.010	252	13255	0.334	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.738	276	15568	0.330	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.758	278	12436	0.331	ng/u1	98
29) Benzo(g,h,i)perylene	27.533	276	13669	0.332	ng/u1	99

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

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