

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023392.D
 Acq On : 24 Dec 2022 03:29
 Operator : CG/JU
 Sample : PB149762BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS762

Quant Time: Dec 24 04:10:20 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 : Thu Dec 22 02:28:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	8150	0.400	ng/u1	0.00
4) Naphthalene-d8	10.810	136	27135	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.641	164	15543	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.386	188	34026	0.400	ng/u1	0.00
17) Chrysene-d12	21.578	240	24374	0.400	ng/u1	0.00
23) Perylene-d12	24.028	264	19188	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.345	96	4199	0.474	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.394	152	12315	0.293	ng/u1	-0.01
18) Fluoranthene-d10	19.416	212	26624	0.302	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.382	88	11293	1.246	ng/u1	94
5) Naphthalene	10.859	128	23250	0.292	ng/u1	99
7) 2-Methylnaphthalene	12.471	142	14926	0.293	ng/u1	100
8) 1-Methylnaphthalene	12.685	142	15997	0.305	ng/u1	100
10) Acenaphthylene	14.358	152	21111	0.306	ng/u1	99
11) Acenaphthene	14.701	153	16769	0.289	ng/u1	99
12) Fluorene	15.686	166	18634	0.286	ng/u1	99
14) Pentachlorophenol	17.040	266	5549	0.588	ng/u1	100
15) Phenanthrene	17.428	178	31251	0.282	ng/u1	100
16) Anthracene	17.521	178	27642	0.301	ng/u1	100
19) Fluoranthene	19.448	202	36002	0.304	ng/u1	99
20) Pyrene	19.811	202	37004	0.310	ng/u1	97
21) Benzo(a)anthracene	21.560	228	31897	0.336	ng/u1	99
22) Chrysene	21.616	228	31049	0.322	ng/u1	99
24) Benzo(b)fluoranthene	23.276	252	26917	0.271	ng/u1	98
25) Benzo(k)fluoranthene	23.326	252	26364	0.283	ng/u1	99
26) Benzo(a)pyrene	23.913	252	22127	0.282	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.584	276	25898	0.282	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.604	278	20461	0.288	ng/u1	98
29) Benzo(g,h,i)perylene	27.369	276	21346	0.286	ng/u1	97

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

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