

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090523\
 Data File : BN027370.D
 Acq On : 05 Sep 2023 11:46
 Operator : MA/JU
 Sample : PB155073BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS073

Quant Time: Sep 06 00:41:22 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 02 06:16:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	3234	0.400	ng/ul	0.00
4) Naphthalene-d8	10.883	136	10570	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.694	164	5857	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.443	188	13769	0.400	ng/ul	0.00
17) Chrysene-d12	21.621	240	11371	0.400	ng/ul	0.01
23) Perylene-d12	24.149	264	9274	0.400	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.418	96	3301	0.767	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	6246	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.464	212	15966	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.452	88	7528	1.729	ng/ul#	87
5) Naphthalene	10.933	128	11010	0.359	ng/ul	99
7) 2-Methylnaphthalene	12.533	142	6446	0.336	ng/ul	100
8) 1-Methylnaphthalene	12.748	142	7106	0.348	ng/ul	97
10) Acenaphthylene	14.421	152	9416	0.351	ng/ul	99
11) Acenaphthene	14.754	153	8194	0.355	ng/ul	99
12) Fluorene	15.744	166	9372	0.363	ng/ul	99
14) Pentachlorophenol	17.076	266	1720	0.641	ng/ul	99
15) Phenanthrene	17.485	178	15522	0.346	ng/ul	99
16) Anthracene	17.582	178	12028	0.337	ng/ul	99
19) Fluoranthene	19.492	202	18242	0.349	ng/ul	98
20) Pyrene	19.859	202	18899	0.355	ng/ul	99
21) Benzo(a)anthracene	21.603	228	13808	0.337	ng/ul	99
22) Chrysene	21.659	228	16878	0.353	ng/ul	100
24) Benzo(b)fluoranthene	23.360	252	15976	0.365	ng/ul	97
25) Benzo(k)fluoranthene	23.413	252	15765	0.368	ng/ul	96
26) Benzo(a)pyrene	24.035	252	12012	0.351	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.831	276	13324	0.351	ng/ul	99
28) Dibenzo(a,h)anthracene	26.861	278	9762	0.336	ng/ul	96
29) Benzo(g,h,i)perylene	27.649	276	13235	0.386	ng/ul	97

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

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