

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020400.D
 Acq On : 24 Jun 2022 16:38
 Operator : CG/JU
 Sample : PB145674BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS674

Quant Time: Jun 25 00:55:11 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.754	152	2784	0.400	ng/u1	0.00
4) Naphthalene-d8	10.535	136	9196	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.382	164	6102	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	14149	0.400	ng/u1	0.00
17) Chrysene-d12	21.312	240	12727	0.400	ng/u1	0.00
23) Perylene-d12	23.595	264	10361	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	2349	0.728	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.130	152	5531	0.367	ng/u1	0.00
18) Fluoranthene-d10	19.151	212	14906	0.355	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	6174	1.875	ng/u1#	82
5) Naphthalene	10.584	128	10002	0.352	ng/u1	100
7) 2-Methylnaphthalene	12.201	142	6622	0.348	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	6998	0.388	ng/u1	100
10) Acenaphthylene	14.099	152	10101	0.349	ng/u1	100
11) Acenaphthene	14.442	153	8019	0.346	ng/u1	99
12) Fluorene	15.427	166	9646	0.352	ng/u1	99
14) Pentachlorophenol	16.787	266	1535	0.538	ng/u1	99
15) Phenanthrene	17.162	178	17213	0.350	ng/u1	99
16) Anthracene	17.255	178	14956	0.352	ng/u1	100
19) Fluoranthene	19.179	202	19792	0.341	ng/u1	99
20) Pyrene	19.546	202	20576	0.341	ng/u1	99
21) Benzo(a)anthracene	21.295	228	16798	0.367	ng/u1	99
22) Chrysene	21.347	228	18875	0.371	ng/u1	100
24) Benzo(b)fluoranthene	22.905	252	17582	0.367	ng/u1	99
25) Benzo(k)fluoranthene	22.949	252	17397	0.359	ng/u1	100
26) Benzo(a)pyrene	23.493	252	15978	0.367	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	25.933	276	16277	0.375	ng/u1#	99
28) Dibenzo(a,h)anthracene	25.943	278	13186	0.377	ng/u1	99
29) Benzo(g,h,i)perylene	26.644	276	13947	0.375	ng/u1	99

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

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