

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020441.D
 Acq On : 25 Jun 2022 18:41
 Operator : CG/JU
 Sample : PB145680BS
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS680

Quant Time: Jun 27 03:43:01 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	7.750	152	2680	0.400	ng/ul	0.00
4) Naphthalene-d8	10.535	136	9162	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.377	164	6172	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.120	188	13876	0.400	ng/ul	0.00
17) Chrysene-d12	21.308	240	11832	0.400	ng/ul	0.00
23) Perylene-d12	23.588	264	9559	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	2254	0.726	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.130	152	5647	0.376	ng/ul	0.00
18) Fluoranthene-d10	19.151	212	14251	0.365	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.235	88	5459	1.722	ng/ul#	81
5) Naphthalene	10.584	128	9456	0.334	ng/ul	99
7) 2-Methylnaphthalene	12.201	142	6316	0.334	ng/ul	99
8) 1-Methylnaphthalene	12.421	142	6740	0.376	ng/ul	99
10) Acenaphthylene	14.095	152	9733	0.333	ng/ul	100
11) Acenaphthene	14.442	153	7686	0.328	ng/ul	98
12) Fluorene	15.427	166	9015	0.325	ng/ul	99
14) Pentachlorophenol	16.786	266	1347	0.481	ng/ul	98
15) Phenanthrene	17.162	178	15714	0.326	ng/ul	100
16) Anthracene	17.255	178	13744	0.330	ng/ul	100
19) Fluoranthene	19.179	202	17728	0.328	ng/ul	100
20) Pyrene	19.541	202	18446	0.329	ng/ul	99
21) Benzo(a)anthracene	21.293	228	14733	0.347	ng/ul	99
22) Chrysene	21.343	228	16637	0.352	ng/ul	99
24) Benzo(b)fluoranthene	22.901	252	15386	0.348	ng/ul	98
25) Benzo(k)fluoranthene	22.945	252	15044	0.337	ng/ul	99
26) Benzo(a)pyrene	23.488	252	14075	0.350	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.928	276	14295	0.357	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.941	278	11565	0.358	ng/ul	99
29) Benzo(g,h,i)perylene	26.639	276	12390	0.361	ng/ul	99

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

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