

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041822\
 Data File : BN019481.D
 Acq On : 18 Apr 2022 16:34
 Operator : CG/JU
 Sample : PB144123BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS123

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/25/2022

Quant Time: Apr 19 03:15:41 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 15:28:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	5131	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	15982	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.460	164	9235	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.204	188	15332	0.400	ng/ul	0.00
17) Chrysene-d12	21.392	240	9170	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	6019	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	4021	0.647	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.217	152	7509	0.350	ng/ul	0.00
18) Fluoranthene-d10	19.234	212	14815	0.396	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.207	88	10637m	1.673	ng/ul	
5) Naphthalene	10.672	128	15126	0.333	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	8897	0.315	ng/ul#	100
8) 1-Methylnaphthalene	12.509	142	9570	0.314	ng/ul#	100
10) Acenaphthylene	14.178	152	12282	0.361	ng/ul	100
11) Acenaphthene	14.520	153	10844	0.323	ng/ul	99
12) Fluorene	15.510	166	11363	0.330	ng/ul	99
14) Pentachlorophenol	16.866	266	1116	0.546	ng/ul	91
15) Phenanthrene	17.242	178	17299	0.331	ng/ul	99
16) Anthracene	17.339	178	12539	0.334	ng/ul	99
19) Fluoranthene	19.262	202	18651	0.366	ng/ul	98
20) Pyrene	19.624	202	19891	0.361	ng/ul	98
21) Benzo(a)anthracene	21.378	228	9448	0.361	ng/ul	100
22) Chrysene	21.428	228	12868	0.378	ng/ul	100
24) Benzo(b)fluoranthene	23.026	252	10615	0.372	ng/ul	94
25) Benzo(k)fluoranthene	23.073	252	12195	0.391	ng/ul	98
26) Benzo(a)pyrene	23.643	252	9925	0.401	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.176	276	9184	0.364	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.196	278	6526	0.389	ng/ul	91
29) Benzo(g,h,i)perylene	26.911	276	10204	0.382	ng/ul	99

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

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