

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081222\
 Data File : BN021212.D
 Acq On : 13 Aug 2022 15:08
 Operator : CG/JU
 Sample : PB146959BS
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS959

Quant Time: Aug 14 23:04:45 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 13 04:02:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	2924	0.400	ng/u1	0.00
4) Naphthalene-d8	10.580	136	10198	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.433	164	5965	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.189	188	11002	0.400	ng/u1	0.00
17) Chrysene-d12	21.395	240	7323	0.400	ng/u1	0.00
23) Perylene-d12	23.722	264	5584	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2602	0.696	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.186	152	5440	0.364	ng/u1	0.00
18) Fluoranthene-d10	19.226	212	10556	0.467	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	6992	1.814	ng/u1	85
5) Naphthalene	10.629	128	10004	0.319	ng/u1	99
7) 2-Methylnaphthalene	12.257	142	6019	0.317	ng/u1	99
8) 1-Methylnaphthalene	12.472	142	6654	0.350	ng/u1	100
10) Acenaphthylene	14.156	152	8830	0.317	ng/u1	99
11) Acenaphthene	14.498	153	7159	0.310	ng/u1	99
12) Fluorene	15.493	166	7761	0.308	ng/u1	98
14) Pentachlorophenol	16.859	266	677	0.399	ng/u1	98
15) Phenanthrene	17.231	178	12128	0.308	ng/u1	100
16) Anthracene	17.328	178	9767	0.320	ng/u1	99
19) Fluoranthene	19.254	202	12944	0.399	ng/u1	98
20) Pyrene	19.621	202	13257	0.396	ng/u1	97
21) Benzo(a)anthracene	21.380	228	7716	0.323	ng/u1	99
22) Chrysene	21.430	228	10016	0.318	ng/u1	100
24) Benzo(b)fluoranthene	23.017	252	8327	0.322	ng/u1	98
25) Benzo(k)fluoranthene	23.064	252	8883	0.326	ng/u1	98
26) Benzo(a)pyrene	23.619	252	7789	0.330	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.132	276	8395	0.298	ng/u1	100
28) Dibenzo(a,h)anthracene	26.152	278	6486	0.292	ng/u1	97
29) Benzo(g,h,i)perylene	26.863	276	7930	0.316	ng/u1	98

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

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