

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020970.D
 Acq On : 03 Aug 2022 09:31
 Operator : CG/JU
 Sample : PB146569BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS569

Quant Time: Aug 03 13:41:10 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 04:19:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	1936	0.400	ng/ul	0.00
4) Naphthalene-d8	10.606	136	6329	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.455	164	3666	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.208	188	7256	0.400	ng/ul	0.00
17) Chrysene-d12	21.416	240	6430	0.400	ng/ul	0.00
23) Perylene-d12	23.760	264	5908	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1551	0.657	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.206	152	3131	0.336	ng/ul	0.00
18) Fluoranthene-d10	19.248	212	6718	0.341	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.240	88	3879	1.526	ng/ul#	85
5) Naphthalene	10.650	128	5691	0.329	ng/ul	99
7) 2-Methylnaphthalene	12.278	142	3454	0.328	ng/ul	100
8) 1-Methylnaphthalene	12.498	142	3751	0.323	ng/ul	100
10) Acenaphthylene	14.178	152	5035	0.336	ng/ul	99
11) Acenaphthene	14.520	153	4101	0.327	ng/ul	98
12) Fluorene	15.515	166	4536	0.314	ng/ul	100
14) Pentachlorophenol	16.879	266	657	0.567	ng/ul	99
15) Phenanthrene	17.251	178	7387	0.300	ng/ul	99
16) Anthracene	17.348	178	5996	0.284	ng/ul	99
19) Fluoranthene	19.281	202	8191	0.283	ng/ul	99
20) Pyrene	19.643	202	8585	0.284	ng/ul	98
21) Benzo(a)anthracene	21.404	228	6428	0.275	ng/ul	99
22) Chrysene	21.457	228	8205	0.284	ng/ul	99
24) Benzo(b)fluoranthene	23.050	252	8195	0.291	ng/ul	99
25) Benzo(k)fluoranthene	23.096	252	8312	0.272	ng/ul	99
26) Benzo(a)pyrene	23.655	252	7468	0.281	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.179	276	8843	0.283	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.199	278	7110	0.278	ng/ul	97
29) Benzo(g,h,i)perylene	26.921	276	8034	0.282	ng/ul	98

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

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