

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081522\
 Data File : BN021231.D
 Acq On : 15 Aug 2022 11:44
 Operator : CG/JU
 Sample : PB146983BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS983

Quant Time: Aug 15 12:11:14 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	7.797	152	2514	0.400	ng/u1	0.00
4) Naphthalene-d8	10.584	136	9032	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.437	164	5158	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.192	188	9094	0.400	ng/u1	0.00
17) Chrysene-d12	21.396	240	5752	0.400	ng/u1	0.00
23) Perylene-d12	23.725	264	4642	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2293	0.713	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.190	152	4537	0.343	ng/u1	0.00
18) Fluoranthene-d10	19.225	212	8422	0.474	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	5644	1.704	ng/u1	86
5) Naphthalene	10.634	128	8581	0.309	ng/u1	100
7) 2-Methylnaphthalene	12.261	142	5053	0.300	ng/u1	99
8) 1-Methylnaphthalene	12.476	142	5576	0.331	ng/u1	99
10) Acenaphthylene	14.155	152	7240	0.300	ng/u1	99
11) Acenaphthene	14.497	153	6046	0.303	ng/u1	99
12) Fluorene	15.492	166	6373	0.292	ng/u1	99
14) Pentachlorophenol	16.862	266	503	0.359	ng/u1	99
15) Phenanthrene	17.230	178	9827	0.302	ng/u1	99
16) Anthracene	17.331	178	7514	0.298	ng/u1	99
19) Fluoranthene	19.257	202	10374	0.408	ng/u1	99
20) Pyrene	19.620	202	10738	0.409	ng/u1	96
21) Benzo(a)anthracene	21.381	228	5631	0.300	ng/u1	99
22) Chrysene	21.433	228	7702	0.312	ng/u1	99
24) Benzo(b)fluoranthene	23.018	252	6885	0.320	ng/u1	88
25) Benzo(k)fluoranthene	23.064	252	7335	0.323	ng/u1	97
26) Benzo(a)pyrene	23.626	252	6609	0.337	ng/u1#	70
27) Indeno(1,2,3-cd)pyrene	26.136	276	6997	0.299	ng/u1#	94
28) Dibenzo(a,h)anthracene	26.153	278	5235	0.283	ng/u1	95
29) Benzo(g,h,i)perylene	26.864	276	6474	0.310	ng/u1	98

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

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