

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082222\
 Data File : BN021315.D
 Acq On : 23 Aug 2022 01:01
 Operator : CG/JU
 Sample : PB147177BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS177

Quant Time: Aug 23 04:05:15 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 18:06:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.097	152	2499	0.400	ng/ul	0.00
4) Naphthalene-d8	10.927	136	8141	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.749	164	4436	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.498	188	9532	0.400	ng/ul	0.00
17) Chrysene-d12	21.685	240	7926	0.400	ng/ul	# 0.00
23) Perylene-d12	24.217	264	6836	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.385	96	1950	0.605	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.511	152	3970	0.322	ng/ul	0.00
18) Fluoranthene-d10	19.525	212	8068	0.330	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.422	88	5069	1.501	ng/ul#	89
5) Naphthalene	10.982	128	7029	0.273	ng/ul	99
7) 2-Methylnaphthalene	12.588	142	4496	0.279	ng/ul	98
8) 1-Methylnaphthalene	12.808	142	4695	0.310	ng/ul	99
10) Acenaphthylene	14.472	152	6032	0.263	ng/ul	99
11) Acenaphthene	14.814	153	4701	0.268	ng/ul	99
12) Fluorene	15.799	166	5253	0.268	ng/ul	99
14) Pentachlorophenol	17.143	266	580	0.551	ng/ul	98
15) Phenanthrene	17.540	178	9332	0.264	ng/ul	100
16) Anthracene	17.633	178	7403	0.261	ng/ul	99
19) Fluoranthene	19.552	202	10294	0.274	ng/ul	98
20) Pyrene	19.915	202	10165	0.270	ng/ul	95
21) Benzo(a)anthracene	21.668	228	7946	0.289	ng/ul	99
22) Chrysene	21.723	228	10129	0.287	ng/ul	99
24) Benzo(b)fluoranthene	23.436	252	9086	0.309	ng/ul	98
25) Benzo(k)fluoranthene	23.489	252	9486	0.300	ng/ul#	97
26) Benzo(a)pyrene	24.103	252	8021	0.290	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.894	276	10004	0.308	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.914	278	8317	0.311	ng/ul	95
29) Benzo(g,h,i)perylene	27.713	276	9559	0.296	ng/ul	97

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

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