

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030113.D
 Acq On : 28 Feb 2024 03:58
 Operator : MA/JU
 Sample : PB159254BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS254

Quant Time: Feb 28 04:26:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	3747	0.400	ng/ul	0.00
4) Naphthalene-d8	10.645	136	10559	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.497	164	4678	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.264	188	8223	0.400	ng/ul	0.00
17) Chrysene-d12	21.464	240	6084	0.400	ng/ul	0.00
23) Perylene-d12	23.826	264	5136	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	3729	0.798	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	5038	0.375	ng/ul	0.00
18) Fluoranthene-d10	19.295	212	8444	0.422	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.349	88	10505	2.065	ng/ul#	83
5) Naphthalene	10.694	128	11602	0.390	ng/ul	99
7) 2-Methylnaphthalene	12.317	142	6203	0.364	ng/ul	98
8) 1-Methylnaphthalene	12.537	142	6573	0.379	ng/ul	99
10) Acenaphthylene	14.220	152	8720	0.382	ng/ul	97
11) Acenaphthene	14.562	153	7074	0.408	ng/ul	98
12) Fluorene	15.566	166	7136	0.393	ng/ul	100
14) Pentachlorophenol	16.913	266	1386	0.756	ng/ul#	49
15) Phenanthrene	17.302	178	10113	0.394	ng/ul	97
16) Anthracene	17.403	178	8489	0.393	ng/ul	96
19) Fluoranthene	19.323	202	11348	0.406	ng/ul	98
20) Pyrene	19.685	202	12434	0.434	ng/ul	99
21) Benzo(a)anthracene	21.449	228	7567	0.345	ng/ul	99
22) Chrysene	21.499	228	10046	0.418	ng/ul	99
24) Benzo(b)fluoranthene	23.104	252	8789	0.455	ng/ul	96
25) Benzo(k)fluoranthene	23.153	252	10335	0.482	ng/ul	97
26) Benzo(a)pyrene	23.721	252	8261	0.455	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.299	276	9355	0.424	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.329	278	7167	0.414	ng/ul#	88
29) Benzo(g,h,i)perylene	27.016	276	8936	0.446	ng/ul	98

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

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