

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020860.D
 Acq On : 26 Jul 2022 13:51
 Operator : CG/JU
 Sample : PB146371BS
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS371

Quant Time: Jul 26 15:14:47 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 18:06:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	2046	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	7352	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	4325	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	9115	0.400	ng/ul	0.00
17) Chrysene-d12	21.426	240	8044	0.400	ng/ul	0.00
23) Perylene-d12	23.770	264	6616	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	2056	0.836	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	5061	0.453	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	11865	0.462	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4913	1.945	ng/ul#	81
5) Naphthalene	10.661	128	8818	0.383	ng/ul	100
7) 2-Methylnaphthalene	12.289	142	5654	0.388	ng/ul	99
8) 1-Methylnaphthalene	12.504	142	6015	0.426	ng/ul	100
10) Acenaphthylene	14.187	152	8255	0.384	ng/ul	99
11) Acenaphthene	14.530	153	6591	0.379	ng/ul	99
12) Fluorene	15.520	166	7426	0.389	ng/ul	99
14) Pentachlorophenol	16.884	266	1308	0.953	ng/ul	98
15) Phenanthrene	17.259	178	12467	0.373	ng/ul	100
16) Anthracene	17.356	178	10350	0.387	ng/ul	100
19) Fluoranthene	19.286	202	14585	0.387	ng/ul	99
20) Pyrene	19.648	202	14971	0.381	ng/ul	99
21) Benzo(a)anthracene	21.412	228	10751	0.405	ng/ul	100
22) Chrysene	21.461	228	13330	0.390	ng/ul	100
24) Benzo(b)fluoranthene	23.060	252	12096	0.386	ng/ul	98
25) Benzo(k)fluoranthene	23.107	252	12774	0.394	ng/ul	99
26) Benzo(a)pyrene	23.668	252	11023	0.381	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.205	276	11720	0.385	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.225	278	9147	0.394	ng/ul	99
29) Benzo(g,h,i)perylene	26.939	276	10611	0.387	ng/ul	99

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

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