

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020857.D
 Acq On : 26 Jul 2022 12:00
 Operator : CG/JU
 Sample : PB146403BS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS403

Quant Time: Jul 26 13:50:49 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	2058	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	7108	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	4360	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	9163	0.400	ng/ul	0.00
17) Chrysene-d12	21.423	240	8047	0.400	ng/ul	0.00
23) Perylene-d12	23.773	264	6641	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1945	0.787	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.212	152	4765	0.441	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	11185	0.435	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4858	1.912	ng/ul#	81
5) Naphthalene	10.661	128	8256	0.371	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	5357	0.380	ng/ul	100
8) 1-Methylnaphthalene	12.503	142	5731	0.420	ng/ul	99
10) Acenaphthylene	14.187	152	8215	0.379	ng/ul	100
11) Acenaphthene	14.530	153	6474	0.369	ng/ul	100
12) Fluorene	15.520	166	7276	0.378	ng/ul	100
14) Pentachlorophenol	16.884	266	1317	0.954	ng/ul	99
15) Phenanthrene	17.259	178	12044	0.359	ng/ul	100
16) Anthracene	17.352	178	10222	0.380	ng/ul	99
19) Fluoranthene	19.285	202	13933	0.370	ng/ul	99
20) Pyrene	19.648	202	14435	0.367	ng/ul	100
21) Benzo(a)anthracene	21.408	228	10629	0.400	ng/ul	99
22) Chrysene	21.461	228	12617	0.369	ng/ul	99
24) Benzo(b)fluoranthene	23.057	252	11958	0.380	ng/ul	96
25) Benzo(k)fluoranthene	23.104	252	11992	0.368	ng/ul	98
26) Benzo(a)pyrene	23.665	252	10861	0.374	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.201	276	11279	0.369	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.221	278	8796	0.378	ng/ul	98
29) Benzo(g,h,i)perylene	26.939	276	10108	0.367	ng/ul	99

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

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