

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020858.D
 Acq On : 26 Jul 2022 12:37
 Operator : CG/JU
 Sample : PB146405BS
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS405

Quant Time: Jul 26 13:51:30 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	2176	0.400	ng/ul	0.00
4) Naphthalene-d8	10.612	136	7740	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.465	164	4556	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	9451	0.400	ng/ul	0.00
17) Chrysene-d12	21.425	240	8313	0.400	ng/ul	0.00
23) Perylene-d12	23.769	264	7049	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1939	0.742	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.212	152	4866	0.414	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	11059	0.416	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4767	1.774	ng/ul#	80
5) Naphthalene	10.661	128	8613	0.355	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	5507	0.359	ng/ul	99
8) 1-Methylnaphthalene	12.504	142	5723	0.385	ng/ul	99
10) Acenaphthylene	14.187	152	8052	0.355	ng/ul	100
11) Acenaphthene	14.530	153	6473	0.353	ng/ul	99
12) Fluorene	15.524	166	7258	0.361	ng/ul	99
14) Pentachlorophenol	16.884	266	1278	0.898	ng/ul	99
15) Phenanthrene	17.259	178	12083	0.349	ng/ul	99
16) Anthracene	17.356	178	10024	0.361	ng/ul	100
19) Fluoranthene	19.285	202	13761	0.353	ng/ul	99
20) Pyrene	19.648	202	14370	0.354	ng/ul	99
21) Benzo(a)anthracene	21.410	228	10307	0.375	ng/ul	100
22) Chrysene	21.463	228	12763	0.361	ng/ul	99
24) Benzo(b)fluoranthene	23.059	252	11780	0.353	ng/ul	98
25) Benzo(k)fluoranthene	23.105	252	12512	0.362	ng/ul	99
26) Benzo(a)pyrene	23.667	252	10938	0.355	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.203	276	11695	0.361	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.217	278	9126	0.369	ng/ul	99
29) Benzo(g,h,i)perylene	26.938	276	10305	0.353	ng/ul	100

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

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