

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020885.D
 Acq On : 27 Jul 2022 05:51
 Operator : CG/JU
 Sample : PB146442BS
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS442

Quant Time: Jul 27 06:21:07 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	7.813	152	2426	0.400	ng/ul	0.00
4) Naphthalene-d8	10.606	136	8393	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.460	164	5286	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	11162	0.400	ng/ul	0.00
17) Chrysene-d12	21.417	240	9941	0.400	ng/ul	-0.01
23) Perylene-d12	23.761	264	8368	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1928	0.661	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.207	152	4938	0.387	ng/ul	-0.01
18) Fluoranthene-d10	19.253	212	11520	0.363	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4724	1.577	ng/ul#	84
5) Naphthalene	10.656	128	8581	0.326	ng/ul	99
7) 2-Methylnaphthalene	12.284	142	5600	0.337	ng/ul	100
8) 1-Methylnaphthalene	12.498	142	5959	0.370	ng/ul	99
10) Acenaphthylene	14.183	152	8585	0.327	ng/ul	100
11) Acenaphthene	14.525	153	6698	0.315	ng/ul	99
12) Fluorene	15.515	166	7554	0.324	ng/ul	98
14) Pentachlorophenol	16.875	266	1728	1.028	ng/ul	99
15) Phenanthrene	17.255	178	12514	0.306	ng/ul	100
16) Anthracene	17.348	178	10758	0.328	ng/ul	99
19) Fluoranthene	19.281	202	14406	0.309	ng/ul	99
20) Pyrene	19.643	202	14950	0.308	ng/ul	100
21) Benzo(a)anthracene	21.403	228	11721	0.357	ng/ul	100
22) Chrysene	21.455	228	13658	0.323	ng/ul	100
24) Benzo(b)fluoranthene	23.051	252	12608	0.318	ng/ul	98
25) Benzo(k)fluoranthene	23.098	252	13076	0.319	ng/ul	98
26) Benzo(a)pyrene	23.659	252	11862	0.324	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.195	276	12248	0.318	ng/ul	97
28) Dibenzo(a,h)anthracene	26.205	278	9734	0.332	ng/ul	98
29) Benzo(g,h,i)perylene	26.932	276	10744	0.310	ng/ul	98

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

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