

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071922\
 Data File : BN020793.D
 Acq On : 20 Jul 2022 03:04
 Operator : CG/JU
 Sample : PB146299BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS299

Quant Time: Jul 20 03:39:27 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 01:02:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	2238	0.400	ng/ul	0.00
4) Naphthalene-d8	10.617	136	8005	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.470	164	3928	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	10602	0.400	ng/ul	0.00
17) Chrysene-d12	21.423	240	10566	0.400	ng/ul	0.00
23) Perylene-d12	23.770	264	8712	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1872	0.722	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	4847	0.369	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	12707	0.365	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	4651	1.757	ng/ul#	82
5) Naphthalene	10.667	128	8777	0.355	ng/ul	100
7) 2-Methylnaphthalene	12.295	142	5595	0.338	ng/ul	100
8) 1-Methylnaphthalene	12.509	142	6045	0.386	ng/ul	100
10) Acenaphthylene	14.192	152	7347	0.395	ng/ul	100
11) Acenaphthene	14.530	153	5445	0.365	ng/ul	99
12) Fluorene	15.529	166	6231	0.353	ng/ul	98
14) Pentachlorophenol	16.888	266	1052	0.492	ng/ul	100
15) Phenanthrene	17.264	178	13792	0.374	ng/ul	99
16) Anthracene	17.361	178	11254	0.354	ng/ul	99
19) Fluoranthene	19.290	202	16199	0.336	ng/ul	99
20) Pyrene	19.653	202	16926	0.338	ng/ul	98
21) Benzo(a)anthracene	21.409	228	12824	0.338	ng/ul	99
22) Chrysene	21.458	228	16669	0.395	ng/ul	100
24) Benzo(b)fluoranthene	23.057	252	15510	0.385	ng/ul	97
25) Benzo(k)fluoranthene	23.104	252	16307	0.401	ng/ul	100
26) Benzo(a)pyrene	23.668	252	13883	0.379	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.198	276	15314	0.420	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.222	278	12048	0.409	ng/ul	99
29) Benzo(g,h,i)perylene	26.939	276	14329	0.458	ng/ul	100

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

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