

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020890.D
 Acq On : 27 Jul 2022 09:28
 Operator : CG/JU
 Sample : PB146444BS
 Misc :
 ALS Vial : 72 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS444

Quant Time: Jul 27 18:02:03 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	7.817	152	2215	0.400	ng/ul	0.00
4) Naphthalene-d8	10.611	136	7773	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.464	164	4860	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.216	188	10208	0.400	ng/ul	0.00
17) Chrysene-d12	21.426	240	8561	0.400	ng/ul	0.00
23) Perylene-d12	23.770	264	6985	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1892	0.711	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.211	152	4820	0.408	ng/ul	0.00
18) Fluoranthene-d10	19.251	212	11163	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.248	88	4617	1.688	ng/ul#	79
5) Naphthalene	10.660	128	8463	0.347	ng/ul	99
7) 2-Methylnaphthalene	12.282	142	5540	0.360	ng/ul	99
8) 1-Methylnaphthalene	12.502	142	5904	0.396	ng/ul	99
10) Acenaphthylene	14.181	152	8278	0.342	ng/ul	99
11) Acenaphthene	14.524	153	6651	0.340	ng/ul	100
12) Fluorene	15.518	166	7466	0.348	ng/ul	98
14) Pentachlorophenol	16.878	266	1385	0.901	ng/ul	97
15) Phenanthrene	17.258	178	12366	0.330	ng/ul	99
16) Anthracene	17.351	178	10303	0.344	ng/ul	99
19) Fluoranthene	19.284	202	14007	0.349	ng/ul	98
20) Pyrene	19.646	202	14448	0.345	ng/ul	100
21) Benzo(a)anthracene	21.411	228	10595	0.375	ng/ul	99
22) Chrysene	21.461	228	12741	0.350	ng/ul	100
24) Benzo(b)fluoranthene	23.057	252	11746	0.355	ng/ul	97
25) Benzo(k)fluoranthene	23.107	252	12003	0.351	ng/ul	99
26) Benzo(a)pyrene	23.668	252	10656	0.349	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.201	276	11056	0.344	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.222	278	8683	0.355	ng/ul	99
29) Benzo(g,h,i)perylene	26.939	276	9989	0.345	ng/ul	99

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

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