

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016086.D
 Acq On : 27 Aug 2021 01:48
 Operator : CG/JU
 Sample : PB138647BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS647

Quant Time: Aug 27 03:31:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 26 11:35:43 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	1433	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	6202	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	3731	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.247	188	8093	0.400	ng/ul	0.00
17) Chrysene-d12	21.429	240	8355	0.400	ng/ul	0.00
23) Perylene-d12	23.817	264	7576	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1365	0.703	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	3539	0.379	ng/ul	0.00
18) Fluoranthene-d10	19.277	212	9107	0.405	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	3321	1.694	ng/ul#	86
5) Naphthalene	10.716	128	5738	0.358	ng/ul	99
7) 2-Methylnaphthalene	12.327	142	3919	0.362	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	3650	0.344	ng/ul	99
10) Acenaphthylene	14.220	152	4829	0.367	ng/ul	99
11) Acenaphthene	14.562	153	3988	0.351	ng/ul	98
12) Fluorene	15.548	166	4648	0.345	ng/ul	97
14) Pentachlorophenol	16.897	266	924	0.545	ng/ul	98
15) Phenanthrene	17.289	178	7911	0.354	ng/ul	99
16) Anthracene	17.378	178	6922	0.359	ng/ul	100
19) Fluoranthene	19.305	202	9591	0.367	ng/ul	99
20) Pyrene	19.667	202	9712	0.370	ng/ul	99
21) Benzo(a)anthracene	21.415	228	9784	0.376	ng/ul	99
22) Chrysene	21.467	228	10792	0.386	ng/ul	100
24) Benzo(b)fluoranthene	23.087	252	10822	0.383	ng/ul	99
25) Benzo(k)fluoranthene	23.136	252	10874	0.382	ng/ul	100
26) Benzo(a)pyrene	23.709	252	9093	0.383	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	26.286	276	12249	0.389	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.302	278	10223	0.386	ng/ul	99
29) Benzo(g,h,i)perylene	27.040	276	10219	0.377	ng/ul	99

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

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