

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN12221\
 Data File : BN018170.D
 Acq On : 23 Dec 2021 02:46
 Operator : CG/JU
 Sample : PB141605BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS605

Quant Time: Dec 23 03:26:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 14:08:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	3076	0.400	ng/ul	0.00
4) Naphthalene-d8	10.394	136	9923	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.249	164	6329	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.995	188	14174	0.400	ng/ul	0.00
17) Chrysene-d12	21.189	240	13852	0.400	ng/ul	0.00
23) Perylene-d12	23.416	264	12616	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	2445	0.769	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.994	152	5546	0.406	ng/ul	0.00
18) Fluoranthene-d10	19.027	212	16197	0.431	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	7196	1.872	ng/ul	90
5) Naphthalene	10.438	128	11385	0.404	ng/ul	100
7) 2-Methylnaphthalene	12.066	142	7156	0.395	ng/ul	99
8) 1-Methylnaphthalene	12.286	142	7386	0.392	ng/ul	100
10) Acenaphthylene	13.972	152	9897	0.395	ng/ul	99
11) Acenaphthene	14.314	153	8762	0.397	ng/ul	99
12) Fluorene	15.304	166	10127	0.396	ng/ul	100
14) Pentachlorophenol	16.662	266	1490	0.559	ng/ul	97
15) Phenanthrene	17.038	178	18488	0.400	ng/ul	99
16) Anthracene	17.130	178	14743	0.394	ng/ul	99
19) Fluoranthene	19.060	202	22941	0.432	ng/ul	100
20) Pyrene	19.422	202	24012	0.431	ng/ul	99
21) Benzo(a)anthracene	21.174	228	17980	0.390	ng/ul	99
22) Chrysene	21.227	228	24227	0.429	ng/ul	100
24) Benzo(b)fluoranthene	22.746	252	21513	0.445	ng/ul	99
25) Benzo(k)fluoranthene	22.790	252	26378	0.438	ng/ul	100
26) Benzo(a)pyrene	23.319	252	19197	0.420	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.674	276	24215	0.402	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.694	278	18230	0.399	ng/ul	98
29) Benzo(g,h,i)perylene	26.365	276	20354	0.409	ng/ul	99

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

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