

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071922\
 Data File : BN020780.D
 Acq On : 19 Jul 2022 18:23
 Operator : CG/JU
 Sample : PB146295BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS295

Quant Time: Jul 20 01:05:29 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	2291	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	9167	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.475	164	5530	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.226	188	11694	0.400	ng/ul	0.00
17) Chrysene-d12	21.430	240	9228	0.400	ng/ul	0.00
23) Perylene-d12	23.780	264	8600	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	2169	0.817	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	5633	0.374	ng/ul	0.00
18) Fluoranthene-d10	19.263	212	13745	0.452	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.256	88	5783	2.134	ng/ul#	84
5) Naphthalene	10.667	128	11010	0.389	ng/ul	100
7) 2-Methylnaphthalene	12.294	142	7072	0.373	ng/ul	99
8) 1-Methylnaphthalene	12.514	142	7512	0.418	ng/ul	100
10) Acenaphthylene	14.193	152	10326	0.394	ng/ul	100
11) Acenaphthene	14.535	153	8427	0.402	ng/ul	100
12) Fluorene	15.530	166	9482	0.381	ng/ul	99
14) Pentachlorophenol	16.888	266	1333	0.565	ng/ul	99
15) Phenanthrene	17.264	178	16246	0.400	ng/ul	99
16) Anthracene	17.361	178	13347	0.380	ng/ul	100
19) Fluoranthene	19.291	202	19233	0.456	ng/ul	98
20) Pyrene	19.653	202	19872	0.455	ng/ul	97
21) Benzo(a)anthracene	21.415	228	12578	0.379	ng/ul	100
22) Chrysene	21.468	228	15531	0.421	ng/ul	100
24) Benzo(b)fluoranthene	23.066	252	15531	0.390	ng/ul	97
25) Benzo(k)fluoranthene	23.113	252	16366	0.407	ng/ul	99
26) Benzo(a)pyrene	23.677	252	15499	0.429	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.212	276	16977	0.472	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.232	278	13136	0.452	ng/ul	100
29) Benzo(g,h,i)perylene	26.950	276	15143	0.490	ng/ul	99

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

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