

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026411.D
 Acq On : 09 Jul 2023 07:30
 Operator : MA/JU
 Sample : PB153886BS
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS886

Quant Time: Jul 09 15:49:36 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 08 09:20:18 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	4149	0.400	ng/ul	0.00
4) Naphthalene-d8	10.736	136	15028	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.570	164	8241	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.330	188	12304	0.400	ng/ul	0.00
17) Chrysene-d12	21.516	240	7840	0.400	ng/ul	-0.01
23) Perylene-d12	23.975	264	7610	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3107	0.665	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.336	152	6010	0.260	ng/ul	0.00
18) Fluoranthene-d10	19.358	212	10060	0.337	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	8698	1.753	ng/ul#	86
5) Naphthalene	10.785	128	15179	0.352	ng/ul	99
7) 2-Methylnaphthalene	12.408	142	8486	0.309	ng/ul	100
8) 1-Methylnaphthalene	12.617	142	9453	0.333	ng/ul	99
10) Acenaphthylene	14.292	152	13849	0.359	ng/ul	98
11) Acenaphthene	14.630	153	11370	0.360	ng/ul	99
12) Fluorene	15.629	166	11116	0.307	ng/ul	99
14) Pentachlorophenol	16.992	266	1692	0.384	ng/ul	99
15) Phenanthrene	17.368	178	14299	0.352	ng/ul	98
16) Anthracene	17.465	178	11936	0.330	ng/ul	96
19) Fluoranthene	19.386	202	15077	0.360	ng/ul	99
20) Pyrene	19.748	202	16393	0.376	ng/ul	97
21) Benzo(a)anthracene	21.502	228	10516	0.325	ng/ul	99
22) Chrysene	21.554	228	12078	0.359	ng/ul	99
24) Benzo(b)fluoranthene	23.223	252	11947	0.368	ng/ul	92
25) Benzo(k)fluoranthene	23.270	252	13305	0.408	ng/ul	94
26) Benzo(a)pyrene	23.870	252	10973	0.389	ng/ul	90
27) Indeno(1,2,3-cd)pyrene	26.547	276	13418	0.435	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.563	278	10069	0.432	ng/ul	93
29) Benzo(g,h,i)perylene	27.338	276	12034	0.439	ng/ul	98

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

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