

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081823\
 Data File : BN026983.D
 Acq On : 19 Aug 2023 06:53
 Operator : MA/JU
 Sample : PB154872BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS872

Quant Time: Aug 19 07:23:34 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 18 13:11:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.101	152	10604	0.400	ng/ul	0.00
4) Naphthalene-d8	10.927	136	33863	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.726	164	22148	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.473	188	48906	0.400	ng/ul	0.00
17) Chrysene-d12	21.647	240	37967	0.400	ng/ul	0.00
23) Perylene-d12	24.190	264	33516	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.439	96	5744	0.456	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.500	152	14690	0.295	ng/ul	0.00
18) Fluoranthene-d10	19.492	212	33868	0.282	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.477	88	13696	1.066	ng/ul#	93
5) Naphthalene	10.977	128	21894	0.229	ng/ul	100
7) 2-Methylnaphthalene	12.572	142	15017	0.259	ng/ul	100
8) 1-Methylnaphthalene	12.792	142	16682	0.275	ng/ul	100
10) Acenaphthylene	14.453	152	23151	0.187	ng/ul	99
11) Acenaphthene	14.791	153	18374	0.214	ng/ul	100
12) Fluorene	15.772	166	20443	0.210	ng/ul	98
14) Pentachlorophenol	17.097	266	3985	0.288	ng/ul	100
15) Phenanthrene	17.515	178	34676	0.216	ng/ul	99
16) Anthracene	17.608	178	29056	0.200	ng/ul	100
19) Fluoranthene	19.520	202	39928	0.253	ng/ul	98
20) Pyrene	19.887	202	40913	0.243	ng/ul	98
21) Benzo(a)anthracene	21.630	228	33940	0.225	ng/ul	99
22) Chrysene	21.688	228	35477	0.238	ng/ul	99
24) Benzo(b)fluoranthene	23.401	252	33462	0.264	ng/ul	97
25) Benzo(k)fluoranthene	23.454	252	32584	0.261	ng/ul#	95
26) Benzo(a)pyrene	24.076	252	26650	0.233	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.881	276	30923	0.207	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.911	278	24193	0.203	ng/ul	98
29) Benzo(g,h,i)perylene	27.719	276	25973	0.208	ng/ul	98

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

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