

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100423\
 Data File : BN028038.D
 Acq On : 04 Oct 2023 16:55
 Operator : MA/JU
 Sample : PB155830BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS830

Quant Time: Oct 05 02:05:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 03 15:50:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	4368	0.400	ng/ul	0.00
4) Naphthalene-d8	10.762	136	13531	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.587	164	7186	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.346	188	15104	0.400	ng/ul	0.00
17) Chrysene-d12	21.530	240	14968	0.400	ng/ul	0.00
23) Perylene-d12	24.000	264	17400	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.326	96	3726	0.679	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	8139	0.373	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	17147	0.361	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.359	88	8858	1.525	ng/ul	92
5) Naphthalene	10.812	128	12872	0.335	ng/ul	100
7) 2-Methylnaphthalene	12.418	142	7986	0.329	ng/ul	100
8) 1-Methylnaphthalene	12.638	142	8374	0.334	ng/ul	100
10) Acenaphthylene	14.314	152	11881	0.364	ng/ul	99
11) Acenaphthene	14.652	153	8640	0.333	ng/ul	100
12) Fluorene	15.637	166	9534	0.330	ng/ul	98
14) Pentachlorophenol	16.978	266	2633	0.813	ng/ul	99
15) Phenanthrene	17.384	178	15226	0.338	ng/ul	99
16) Anthracene	17.477	178	13881	0.345	ng/ul	99
19) Fluoranthene	19.399	202	18197	0.318	ng/ul	98
20) Pyrene	19.766	202	19338	0.326	ng/ul	99
21) Benzo(a)anthracene	21.513	228	19287	0.353	ng/ul	99
22) Chrysene	21.568	228	19033	0.345	ng/ul	99
24) Benzo(b)fluoranthene	23.234	252	21596	0.335	ng/ul	94
25) Benzo(k)fluoranthene	23.281	252	21584	0.336	ng/ul#	94
26) Benzo(a)pyrene	23.889	252	18998	0.345	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.596	276	26447	0.397	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.616	278	21592	0.399	ng/ul	96
29) Benzo(g,h,i)perylene	27.407	276	22219	0.408	ng/ul	96

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

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