

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122623\
 Data File : BM043625.D
 Acq On : 28 Dec 2023 01:37
 Operator : MA/JU
 Sample : PB158076BS
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS076

Quant Time: Dec 28 02:11:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 22:58:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	5752	0.400	ng/ul	0.00
4) Naphthalene-d8	10.774	136	17036	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	7013	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	13514	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.527	240	8008	0.400	ng/ul	# 0.00
23) Perylene-d12	23.942	264	9011	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.377	96	4889	0.673	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	6449	0.267	ng/ul	0.00
18) Fluoranthene-d10	19.376	212	8215	0.287	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.410	88	14406	1.976	ng/ul#	82
5) Naphthalene	10.823	128	14558	0.284	ng/ul	99
7) 2-Methylnaphthalene	12.440	142	8369	0.253	ng/ul	97
8) 1-Methylnaphthalene	12.654	142	8126	0.245	ng/ul	99
10) Acenaphthylene	14.328	152	11323	0.293	ng/ul	98
11) Acenaphthene	14.666	153	8290	0.295	ng/ul	99
12) Fluorene	15.656	166	8908	0.280	ng/ul	98
14) Pentachlorophenol	16.995	266	1854	0.619	ng/ul	97
15) Phenanthrene	17.392	178	13240	0.294	ng/ul	98
16) Anthracene	17.485	178	12197	0.298	ng/ul	99
19) Fluoranthene	19.404	202	12810	0.284	ng/ul#	91
20) Pyrene	19.766	202	13079	0.292	ng/ul#	94
21) Benzo(a)anthracene	21.513	228	10124	0.288	ng/ul	99
22) Chrysene	21.562	228	11690	0.315	ng/ul	99
24) Benzo(b)fluoranthene	23.205	252	11286	0.276	ng/ul	90
25) Benzo(k)fluoranthene	23.255	252	13016	0.307	ng/ul#	88
26) Benzo(a)pyrene	23.836	252	10440	0.309	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.455	276	14230	0.370	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.485	278	11490	0.377	ng/ul#	92
29) Benzo(g,h,i)perylene	27.220	276	12227	0.373	ng/ul	93

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

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