

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101023\
 Data File : BM042265.D
 Acq On : 11 Oct 2023 00:24
 Operator : MA/JU
 Sample : PB156094BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS094

Quant Time: Oct 11 00:59:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 10 14:01:53 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.026	152	3658	0.400	ng/ul	0.00
4) Naphthalene-d8	10.845	136	9018	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.661	164	3866	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.413	188	7178	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.591	240	2887	0.400	ng/ul	# 0.00
23) Perylene-d12	24.064	264	2907	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	3880	0.796	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.423	152	4432	0.358	ng/ul	0.00
18) Fluoranthene-d10	19.436	212	4977	0.459	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.414	88	11392	2.239	ng/ul#	76
5) Naphthalene	10.895	128	11202	0.372	ng/ul	100
7) 2-Methylnaphthalene	12.495	142	6629	0.368	ng/ul	97
8) 1-Methylnaphthalene	12.715	142	6477	0.357	ng/ul	99
10) Acenaphthylene	14.388	152	9364	0.383	ng/ul	99
11) Acenaphthene	14.726	153	7335	0.397	ng/ul	100
12) Fluorene	15.712	166	7981	0.383	ng/ul	98
14) Pentachlorophenol	17.046	266	1772	0.545	ng/ul	99
15) Phenanthrene	17.455	178	11946	0.404	ng/ul	100
16) Anthracene	17.548	178	10133	0.374	ng/ul	99
19) Fluoranthene	19.464	202	11726	0.535	ng/ul#	92
20) Pyrene	19.831	202	11829	0.507	ng/ul#	92
21) Benzo(a)anthracene	21.574	228	8425	0.452	ng/ul	99
22) Chrysene	21.629	228	8770	0.483	ng/ul	99
24) Benzo(b)fluoranthene	23.298	252	8721	0.509	ng/ul	97
25) Benzo(k)fluoranthene	23.348	252	8568	0.508	ng/ul	96
26) Benzo(a)pyrene	23.950	252	7453	0.504	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.656	276	9942	0.475	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.679	278	7317	0.463	ng/ul#	91
29) Benzo(g,h,i)perylene	27.471	276	8589	0.491	ng/ul	94

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

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