

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072623\
 Data File : BM041144.D
 Acq On : 26 Jul 2023 20:27
 Operator : MA/JU
 Sample : PB154169BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS169

Quant Time: Jul 27 01:19:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM072423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 04:20:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	2419	0.400	ng/ul	0.00
4) Naphthalene-d8	10.647	136	7405	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.490	164	3154	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.261	188	5095	0.400	ng/ul	0.00
17) Chrysene-d12	21.454	240	4086	0.400	ng/ul	# 0.00
23) Perylene-d12	23.839	264	3936	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.250	96	2598	0.696	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.247	152	3462	0.327	ng/ul	0.01
18) Fluoranthene-d10	19.292	212	4660	0.419	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.284	88	5824	1.527	ng/ul#	46
5) Naphthalene	10.697	128	6165	0.256	ng/ul	96
7) 2-Methylnaphthalene	12.330	142	3843	0.326	ng/ul	99
8) 1-Methylnaphthalene	12.533	142	4298	0.354	ng/ul	100
10) Acenaphthylene	14.217	152	3812	0.286	ng/ul	97
11) Acenaphthene	14.555	153	3450	0.307	ng/ul	98
12) Fluorene	15.559	166	3316	0.270	ng/ul	98
14) Pentachlorophenol	16.923	266	515	0.347	ng/ul	96
15) Phenanthrene	17.299	178	4698	0.299	ng/ul	97
16) Anthracene	17.401	178	3382	0.266	ng/ul	93
19) Fluoranthene	19.320	202	4983	0.340	ng/ul#	91
20) Pyrene	19.682	202	5401	0.334	ng/ul#	90
21) Benzo(a)anthracene	21.440	228	3689	0.270	ng/ul	97
22) Chrysene	21.489	228	4803	0.304	ng/ul	97
24) Benzo(b)fluoranthene	23.108	252	5050	0.322	ng/ul	90
25) Benzo(k)fluoranthene	23.161	252	4803	0.321	ng/ul#	87
26) Benzo(a)pyrene	23.737	252	4014	0.316	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.307	276	7928	0.422	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.327	278	6201	0.420	ng/ul#	89
29) Benzo(g,h,i)perylene	27.062	276	5456	0.325	ng/ul#	91

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

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