

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072623\
 Data File : BM041143.D
 Acq On : 26 Jul 2023 19:50
 Operator : MA/JU
 Sample : PB154143BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS143

Quant Time: Jul 27 01:19:37 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	2034	0.400	ng/ul	0.00
4) Naphthalene-d8	10.642	136	7253	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.490	164	4124	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	4990	0.400	ng/ul #	0.00
17) Chrysene-d12	21.451	240	3918	0.400	ng/ul #	0.00
23) Perylene-d12	23.839	264	3669	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.254	96	2501	0.797	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.247	152	3288	0.318	ng/ul	0.01
18) Fluoranthene-d10	19.292	212	4350	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.284	88	6596	2.057	ng/ul#	45
5) Naphthalene	10.691	128	7151	0.303	ng/ul	99
7) 2-Methylnaphthalene	12.319	142	3751	0.324	ng/ul	98
8) 1-Methylnaphthalene	12.533	142	4045	0.341	ng/ul	99
10) Acenaphthylene	14.217	152	5803	0.333	ng/ul	99
11) Acenaphthene	14.555	153	5177	0.352	ng/ul	99
12) Fluorene	15.559	166	5193	0.324	ng/ul	98
14) Pentachlorophenol	16.923	266	997	0.686	ng/ul	98
15) Phenanthrene	17.299	178	5335	0.347	ng/ul	97
16) Anthracene	17.401	178	3996	0.321	ng/ul	95
19) Fluoranthene	19.320	202	5628	0.401	ng/ul#	93
20) Pyrene	19.682	202	6097	0.394	ng/ul#	90
21) Benzo(a)anthracene	21.440	228	4239	0.324	ng/ul	97
22) Chrysene	21.489	228	5401	0.357	ng/ul	98
24) Benzo(b)fluoranthene	23.108	252	5471	0.374	ng/ul	93
25) Benzo(k)fluoranthene	23.158	252	5458	0.391	ng/ul#	90
26) Benzo(a)pyrene	23.737	252	4539	0.383	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.307	276	6419	0.366	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.324	278	5063	0.368	ng/ul#	90
29) Benzo(g,h,i)perylene	27.055	276	6038	0.385	ng/ul#	92

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

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