

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072623\
 Data File : BM041142.D
 Acq On : 26 Jul 2023 19:12
 Operator : MA/JU
 Sample : PB154145BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS145

Quant Time: Jul 27 01:19:26 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	7.836	152	3329	0.400	ng/ul	0.00
4) Naphthalene-d8	10.642	136	10774	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.490	164	4747	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	7996	0.400	ng/ul	0.00
17) Chrysene-d12	21.451	240	6030	0.400	ng/ul	# 0.00
23) Perylene-d12	23.836	264	5649	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.250	96	2983	0.581	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.242	152	4430	0.288	ng/ul	0.00
18) Fluoranthene-d10	19.287	212	5992	0.365	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.284	88	8271	1.576	ng/ul#	44
5) Naphthalene	10.691	128	9487	0.271	ng/ul	100
7) 2-Methylnaphthalene	12.319	142	5097	0.297	ng/ul	99
8) 1-Methylnaphthalene	12.528	142	5474	0.310	ng/ul	99
10) Acenaphthylene	14.213	152	7213	0.360	ng/ul	99
11) Acenaphthene	14.555	153	5441	0.322	ng/ul	98
12) Fluorene	15.554	166	5330	0.289	ng/ul	98
14) Pentachlorophenol	16.919	266	974	0.419	ng/ul	97
15) Phenanthrene	17.295	178	7451	0.302	ng/ul	98
16) Anthracene	17.396	178	5718	0.287	ng/ul	96
19) Fluoranthene	19.320	202	7762	0.359	ng/ul#	95
20) Pyrene	19.682	202	11464	0.481	ng/ul	98
21) Benzo(a)anthracene	21.436	228	5990	0.297	ng/ul	99
22) Chrysene	21.486	228	7180	0.308	ng/ul	98
24) Benzo(b)fluoranthene	23.105	252	7384	0.328	ng/ul	96
25) Benzo(k)fluoranthene	23.155	252	7278	0.339	ng/ul#	94
26) Benzo(a)pyrene	23.731	252	6088	0.334	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.300	276	8379	0.310	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.320	278	6613	0.312	ng/ul#	95
29) Benzo(g,h,i)perylene	27.051	276	11068	0.459	ng/ul	97

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

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