

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030914.D
 Acq On : 09 Jul 2021 09:40
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4719

Quant Time: Jul 09 10:17:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 10:44:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.767	152	2025	0.400	ng/ul	0.00
4) Naphthalene-d8	10.553	136	7642	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.402	164	3579	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.151	188	5598	0.400	ng/ul	0.00
17) Chrysene-d12	21.316	240	4419	0.400	ng/ul	0.00
23) Perylene-d12	23.563	264	4285	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.272	96	901	0.408	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.159	152	4585	0.382	ng/ul	0.00
18) Fluoranthene-d10	19.168	212	5646	0.446	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.303	88	854	0.375	ng/ul	96
5) Naphthalene	10.603	128	8730	0.401	ng/ul	98
7) 2-Methylnaphthalene	12.231	142	5497	0.375	ng/ul	98
8) 1-Methylnaphthalene	12.442	142	5206	0.365	ng/ul	99
10) Acenaphthylene	14.125	152	6211	0.389	ng/ul	98
11) Acenaphthene	14.466	153	5372	0.415	ng/ul	99
12) Fluorene	15.463	166	5271	0.369	ng/ul	98
14) Pentachlorophenol	16.825	266	577	0.316	ng/ul#	85
15) Phenanthrene	17.193	178	7052	0.375	ng/ul	97
16) Anthracene	17.290	178	6167	0.374	ng/ul	96
19) Fluoranthene	19.195	202	7746	0.397	ng/ul	98
20) Pyrene	19.557	202	8077	0.412	ng/ul	99
21) Benzo(a)anthracene	21.301	228	6051	0.344	ng/ul	99
22) Chrysene	21.350	228	7786	0.395	ng/ul	98
24) Benzo(b)fluoranthene	22.885	252	7158	0.377	ng/ul	91
25) Benzo(k)fluoranthene	22.929	252	8374	0.428	ng/ul#	90
26) Benzo(a)pyrene	23.466	252	6641	0.411	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	25.851	276	8846	0.421	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.870	278	7121	0.422	ng/ul#	87
29) Benzo(g,h,i)perylene	26.534	276	7917	0.433	ng/ul	97

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

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