

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070821\
 Data File : BM030905.D
 Acq On : 09 Jul 2021 04:11
 Operator : CG/JU
 Sample : PB137515BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS515

Quant Time: Jul 09 05:22:01 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	2365	0.400	ng/ul	0.00
4) Naphthalene-d8	10.554	136	9013	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.398	164	4439	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.145	188	7013	0.400	ng/ul	0.00
17) Chrysene-d12	21.311	240	5200	0.400	ng/ul	0.00
23) Perylene-d12	23.559	264	4983	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	2115	0.821	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.151	152	5377	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.165	212	7356	0.493	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.296	88	5139	1.930	ng/ul#	84
5) Naphthalene	10.604	128	9662	0.376	ng/ul	98
7) 2-Methylnaphthalene	12.227	142	6201	0.359	ng/ul	97
8) 1-Methylnaphthalene	12.443	142	5714	0.340	ng/ul	100
10) Acenaphthylene	14.121	152	7264	0.367	ng/ul	96
11) Acenaphthene	14.463	153	6240	0.389	ng/ul	98
12) Fluorene	15.460	166	6293	0.355	ng/ul	97
14) Pentachlorophenol	16.812	266	1632	0.713	ng/ul	93
15) Phenanthrene	17.186	178	8653	0.367	ng/ul	97
16) Anthracene	17.284	178	7783	0.377	ng/ul	97
19) Fluoranthene	19.195	202	9601	0.418	ng/ul	100
20) Pyrene	19.554	202	9996	0.433	ng/ul	99
21) Benzo(a)anthracene	21.299	228	7555	0.364	ng/ul	99
22) Chrysene	21.345	228	9467	0.408	ng/ul	99
24) Benzo(b)fluoranthene	22.878	252	9410	0.426	ng/ul	93
25) Benzo(k)fluoranthene	22.924	252	10367	0.456	ng/ul	93
26) Benzo(a)pyrene	23.464	252	8211	0.437	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.841	276	10922	0.447	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.856	278	8802	0.449	ng/ul#	91
29) Benzo(g,h,i)perylene	26.527	276	9735	0.458	ng/ul	97

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

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