

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041522\
 Data File : BN019468.D
 Acq On : 15 Apr 2022 18:26
 Operator : CG/JU
 Sample : PB144123BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS123

Quant Time: Apr 18 05:04:04 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 14:35:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	3803	0.400	ng/ul	0.00
4) Naphthalene-d8	10.622	136	12234	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.460	164	7477	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.204	188	13634	0.400	ng/ul	0.00
17) Chrysene-d12	21.392	240	9606	0.400	ng/ul	0.00
23) Perylene-d12	23.745	264	7024	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	2539	0.519	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	4666	0.270	ng/ul	0.00
18) Fluoranthene-d10	19.234	212	23795	0.669	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.211	88	4944	1.049	ng/ul#	77
5) Naphthalene	10.672	128	7098	0.204	ng/ul	98
7) 2-Methylnaphthalene	12.294	142	4169	0.184	ng/ul#	100
8) 1-Methylnaphthalene	12.509	142	4577	0.192	ng/ul#	100
10) Acenaphthylene	14.182	152	6378	0.235	ng/ul	99
11) Acenaphthene	14.520	153	5391	0.199	ng/ul	98
12) Fluorene	15.510	166	5794	0.192	ng/ul	97
14) Pentachlorophenol	16.866	266	2146	0.981	ng/ul	90
15) Phenanthrene	17.242	178	12214	0.270	ng/ul	99
16) Anthracene	17.339	178	10057	0.289	ng/ul	99
19) Fluoranthene	19.262	202	28120	0.588	ng/ul	99
20) Pyrene	19.624	202	30139	0.593	ng/ul	100
21) Benzo(a)anthracene	21.378	228	20446	0.691	ng/ul	100
22) Chrysene	21.430	228	26526	0.685	ng/ul	100
24) Benzo(b)fluoranthene	23.029	252	21480	0.661	ng/ul	93
25) Benzo(k)fluoranthene	23.076	252	25023	0.667	ng/ul	98
26) Benzo(a)pyrene	23.640	252	20270	0.730	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.176	276	19899	0.644	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.189	278	14858	0.676	ng/ul#	90
29) Benzo(g,h,i)perylene	26.910	276	19532	0.654	ng/ul	99

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

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