

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020122\
 Data File : BN018512.D
 Acq On : 01 Feb 2022 13:54
 Operator : CG/JU
 Sample : PB142373BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS373

Quant Time: Feb 02 04:43:45 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 03:59:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	347	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1172	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	611	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.298	188	1162	0.400	ng/ul #	0.00
17) Chrysene-d12	21.468	240	1068	0.400	ng/ul #	0.00
23) Perylene-d12	23.867	264	943	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.295	96	280	0.701	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	599	0.379	ng/ul	0.00
18) Fluoranthene-d10	19.319	212	1229	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.333	88	749	1.783	ng/ul#	72
5) Naphthalene	10.783	128	1195	0.370	ng/ul#	88
7) 2-Methylnaphthalene	12.394	142	709	0.346	ng/ul	100
8) 1-Methylnaphthalene	12.614	142	731	0.349	ng/ul	100
10) Acenaphthylene	14.281	152	961	0.378	ng/ul#	87
11) Acenaphthene	14.623	153	759	0.359	ng/ul	97
12) Fluorene	15.604	166	798	0.353	ng/ul#	98
14) Pentachlorophenol	16.952	266	49	0.475	ng/ul	92
15) Phenanthrene	17.340	178	1382	0.369	ng/ul#	90
16) Anthracene	17.429	178	1097	0.355	ng/ul#	89
19) Fluoranthene	19.351	202	1627	0.394	ng/ul#	84
20) Pyrene	19.709	202	1684	0.397	ng/ul#	82
21) Benzo(a)anthracene	21.453	228	1325	0.377	ng/ul	94
22) Chrysene	21.506	228	1701	0.420	ng/ul	95
24) Benzo(b)fluoranthene	23.137	252	1625	0.413	ng/ul	88
25) Benzo(k)fluoranthene	23.183	252	1680	0.408	ng/ul#	88
26) Benzo(a)pyrene	23.759	252	1436	0.439	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.356	276	1738	0.408	ng/ul#	78
28) Dibenzo(a,h)anthracene	26.376	278	1420	0.424	ng/ul#	77
29) Benzo(g,h,i)perylene	27.118	276	1565	0.422	ng/ul#	78

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

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