

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020122\
 Data File : BN018535.D
 Acq On : 02 Feb 2022 04:07
 Operator : CG/JU
 Sample : PB142347BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS347

Quant Time: Feb 02 05:31:48 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	7.927	152	384	0.400	ng/ul	0.00
4) Naphthalene-d8	10.733	136	1345	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.558	164	738	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.293	188	1470	0.400	ng/ul #	0.00
17) Chrysene-d12	21.464	240	1357	0.400	ng/ul #	0.00
23) Perylene-d12	23.858	264	1243	0.400	ng/ul #-	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	311	0.703	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.322	152	723	0.399	ng/ul	0.00
18) Fluoranthene-d10	19.318	212	1615	0.426	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.332	88	838	1.802	ng/ul#	71
5) Naphthalene	10.783	128	1435	0.387	ng/ul#	91
7) 2-Methylnaphthalene	12.394	142	895	0.381	ng/ul	99
8) 1-Methylnaphthalene	12.608	142	911	0.379	ng/ul	100
10) Acenaphthylene	14.276	152	1306	0.426	ng/ul#	91
11) Acenaphthene	14.618	153	991	0.388	ng/ul	98
12) Fluorene	15.603	166	1095	0.401	ng/ul#	98
14) Pentachlorophenol	16.943	266	311	2.381	ng/ul	100
15) Phenanthrene	17.336	178	1868	0.394	ng/ul	94
16) Anthracene	17.424	178	1566	0.400	ng/ul#	92
19) Fluoranthene	19.346	202	2206	0.421	ng/ul#	87
20) Pyrene	19.709	202	2282	0.424	ng/ul#	88
21) Benzo(a)anthracene	21.447	228	2044	0.458	ng/ul	96
22) Chrysene	21.499	228	2330	0.453	ng/ul	97
24) Benzo(b)fluoranthene	23.130	252	2358	0.455	ng/ul	90
25) Benzo(k)fluoranthene	23.177	252	2296	0.423	ng/ul#	90
26) Benzo(a)pyrene	23.753	252	2164	0.502	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.349	276	2468	0.439	ng/ul#	81
28) Dibenzo(a,h)anthracene	26.366	278	1995	0.452	ng/ul#	82
29) Benzo(g,h,i)perylene	27.107	276	2219	0.454	ng/ul#	82

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

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