

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052422\
 Data File : BN019951.D
 Acq On : 25 May 2022 14:56
 Operator : CG/JU
 Sample : PB145007BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS007

Quant Time: May 26 01:15:27 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN052322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 23 14:58:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.855	152	3302	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	11622	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.470	164	6924	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.214	188	13923	0.400	ng/ul	0.00
17) Chrysene-d12	21.398	240	9438	0.400	ng/ul	0.00
23) Perylene-d12	23.739	264	7538	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2303	0.748	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	7398	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.240	212	14210	0.421	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.349	88	5804	1.806	ng/ul#	71
5) Naphthalene	10.689	128	12688	0.348	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	8442	0.350	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	8824	0.392	ng/ul#	100
10) Acenaphthylene	14.193	152	12832	0.325	ng/ul	99
11) Acenaphthene	14.535	153	9271	0.347	ng/ul	99
12) Fluorene	15.521	166	10366	0.347	ng/ul	99
14) Pentachlorophenol	16.868	266	1349	0.582	ng/ul	90
15) Phenanthrene	17.252	178	16995	0.347	ng/ul	99
16) Anthracene	17.345	178	15029	0.327	ng/ul	99
19) Fluoranthene	19.272	202	18237	0.387	ng/ul	97
20) Pyrene	19.630	202	18107	0.379	ng/ul	99
21) Benzo(a)anthracene	21.383	228	13258	0.344	ng/ul	100
22) Chrysene	21.433	228	14303	0.367	ng/ul	99
24) Benzo(b)fluoranthene	23.029	252	12611	0.378	ng/ul	92
25) Benzo(k)fluoranthene	23.075	252	12445	0.384	ng/ul	98
26) Benzo(a)pyrene	23.634	252	11835	0.376	ng/ul#	79
27) Indeno(1,2,3-cd)pyrene	26.142	276	12799	0.387	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.155	278	10335	0.381	ng/ul#	82
29) Benzo(g,h,i)perylene	26.880	276	10843	0.383	ng/ul	98

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

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