

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071723\
 Data File : BN026539.D
 Acq On : 17 Jul 2023 16:31
 Operator : MA/JU
 Sample : PB153904BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS904

Quant Time: Jul 18 03:05:04 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 17 17:14:21 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.968	152	1830	0.400	ng/ul	0.02
4) Naphthalene-d8	10.765	136	6017	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.584	164	3868	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.334	188	7692	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.520	240	6490	0.400	ng/ul	# 0.00
23) Perylene-d12	23.938	264	6328	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	1832	0.760	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.354	152	2725	0.333	ng/ul	0.00
18) Fluoranthene-d10	19.359	212	7737	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	4530	1.986	ng/ul	90
5) Naphthalene	10.809	128	5188	0.314	ng/ul	97
7) 2-Methylnaphthalene	12.431	142	2992	0.319	ng/ul	99
8) 1-Methylnaphthalene	12.640	142	3378	0.336	ng/ul	100
10) Acenaphthylene	14.311	152	5778	0.305	ng/ul	100
11) Acenaphthene	14.644	153	4510	0.313	ng/ul	99
12) Fluorene	15.639	166	4856	0.321	ng/ul	99
14) Pentachlorophenol	17.001	266	1139	0.587	ng/ul	97
15) Phenanthrene	17.372	178	7370	0.301	ng/ul	98
16) Anthracene	17.474	178	6636	0.312	ng/ul	100
19) Fluoranthene	19.391	202	9460	0.311	ng/ul#	95
20) Pyrene	19.754	202	10163	0.308	ng/ul#	95
21) Benzo(a)anthracene	21.506	228	7345	0.308	ng/ul	98
22) Chrysene	21.558	228	8605	0.317	ng/ul	99
24) Benzo(b)fluoranthene	23.201	252	7487	0.304	ng/ul	96
25) Benzo(k)fluoranthene	23.248	252	8735	0.338	ng/ul#	96
26) Benzo(a)pyrene	23.832	252	7215	0.329	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.457	276	9982	0.347	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.474	278	7995	0.362	ng/ul#	93
29) Benzo(g,h,i)perylene	27.235	276	8739	0.344	ng/ul#	89

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

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