

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030147.D
 Acq On : 29 Feb 2024 01:39
 Operator : MA/JU
 Sample : PB159153BS
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS153

Quant Time: Feb 29 02:39:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	3989	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	11750	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.511	164	5469	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.268	188	9129	0.400	ng/ul	0.00
17) Chrysene-d12	21.466	240	6100	0.400	ng/ul	0.00
23) Perylene-d12	23.833	264	5281	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	4246	0.854	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	6338	0.424	ng/ul	0.01
18) Fluoranthene-d10	19.299	212	9672	0.482	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.353	88	13658	2.522	ng/ul#	81
5) Naphthalene	10.711	128	14168	0.428	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	7991	0.421	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	7960	0.412	ng/ul	99
10) Acenaphthylene	14.229	152	11087	0.415	ng/ul	98
11) Acenaphthene	14.571	153	8786	0.433	ng/ul	99
12) Fluorene	15.570	166	8747	0.412	ng/ul	97
14) Pentachlorophenol	16.917	266	2021	0.993	ng/ul#	49
15) Phenanthrene	17.310	178	12163	0.426	ng/ul	98
16) Anthracene	17.407	178	10297	0.429	ng/ul	98
19) Fluoranthene	19.327	202	13209	0.471	ng/ul	99
20) Pyrene	19.690	202	13819	0.481	ng/ul	99
21) Benzo(a)anthracene	21.454	228	9026	0.410	ng/ul	99
22) Chrysene	21.504	228	10840	0.450	ng/ul	100
24) Benzo(b)fluoranthene	23.111	252	9656	0.486	ng/ul	99
25) Benzo(k)fluoranthene	23.164	252	11849	0.538	ng/ul	98
26) Benzo(a)pyrene	23.734	252	9276	0.496	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.310	276	11014	0.486	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.354	278	8479	0.477	ng/ul#	90
29) Benzo(g,h,i)perylene	27.045	276	10189	0.495	ng/ul	96

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

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